

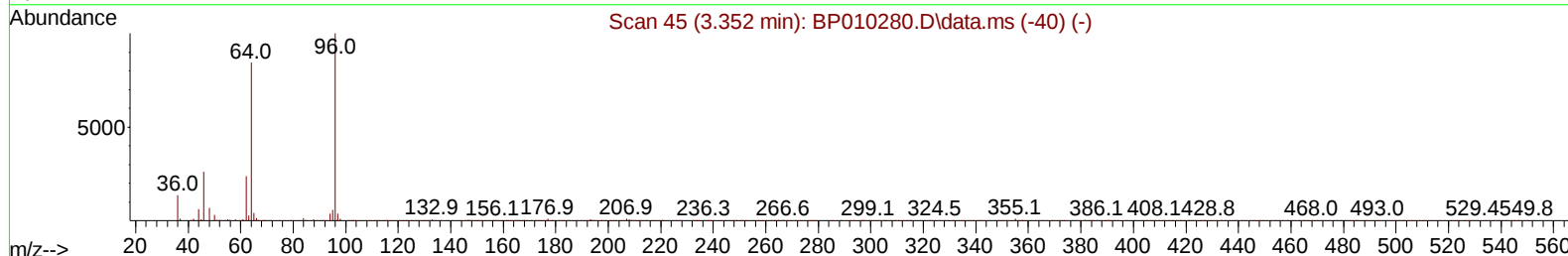
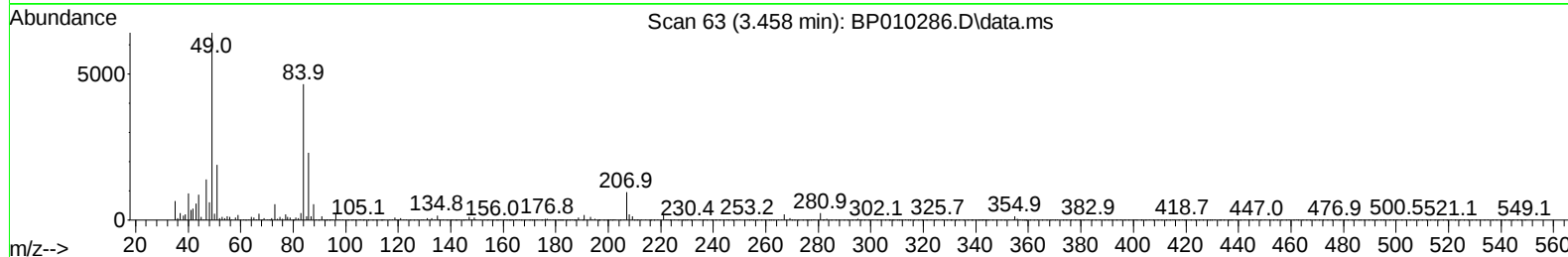
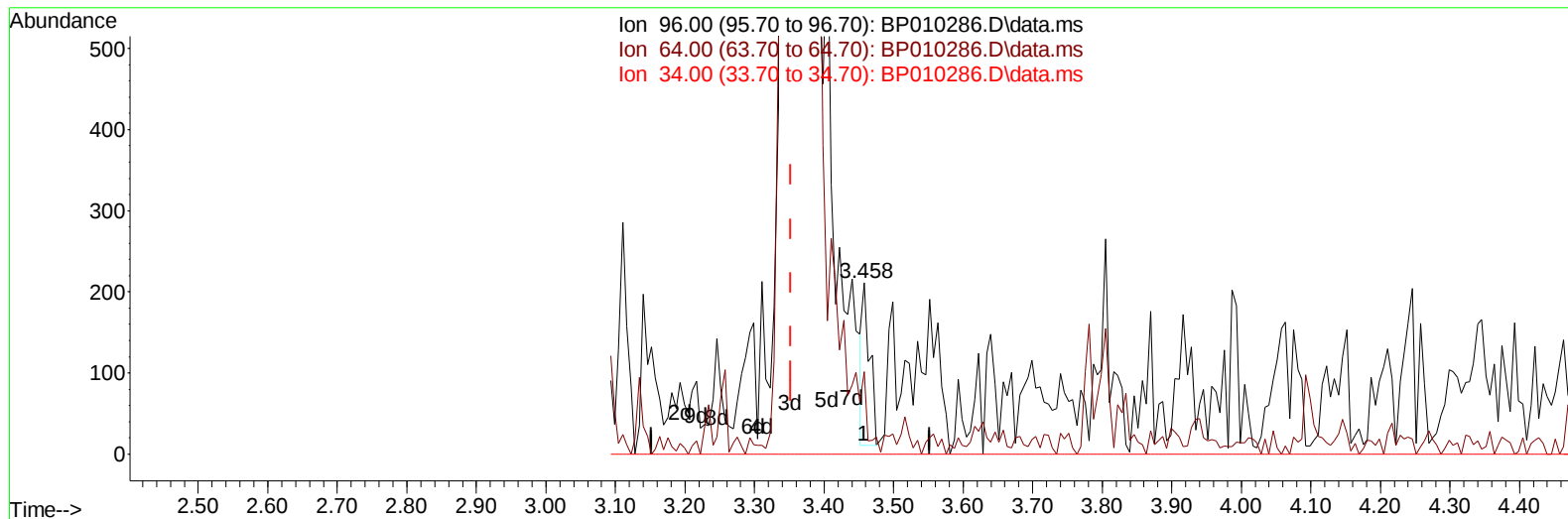
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
 Data File : BP010286.D
 Acq On : 11 May 2022 13:09
 Operator : CG/JU
 Sample : N2632-15
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EW4B7

Manual IntegrationsAPPROVED

Quant Time: May 12 01:12:44 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu May 12 01:10:15 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/12/2022
 Supervised By :Yogesh Patel 05/17/2022



TIC: BP010286.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.458min (+ 0.106) 0.05 ng/uL

response 146

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	48.34
34.00	0.00	0.00
0.00	0.00	0.00

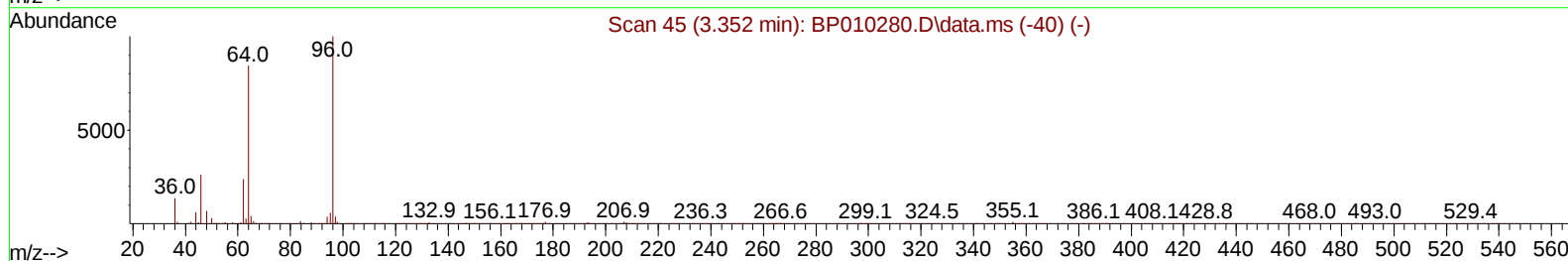
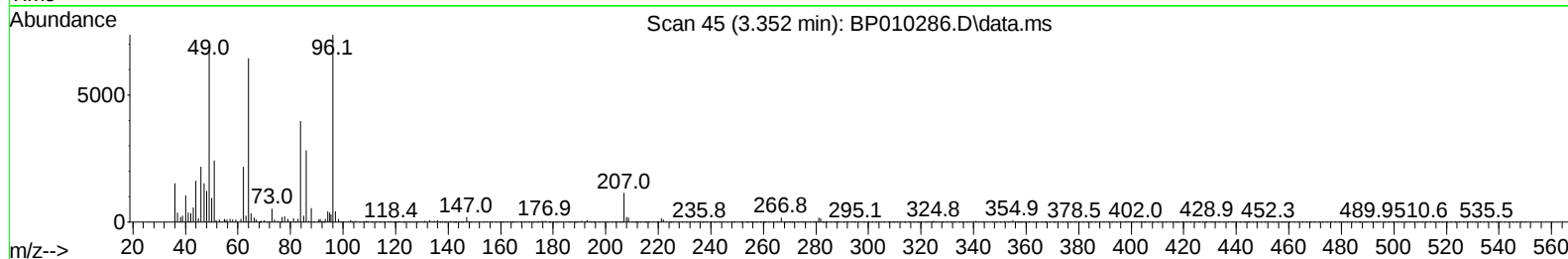
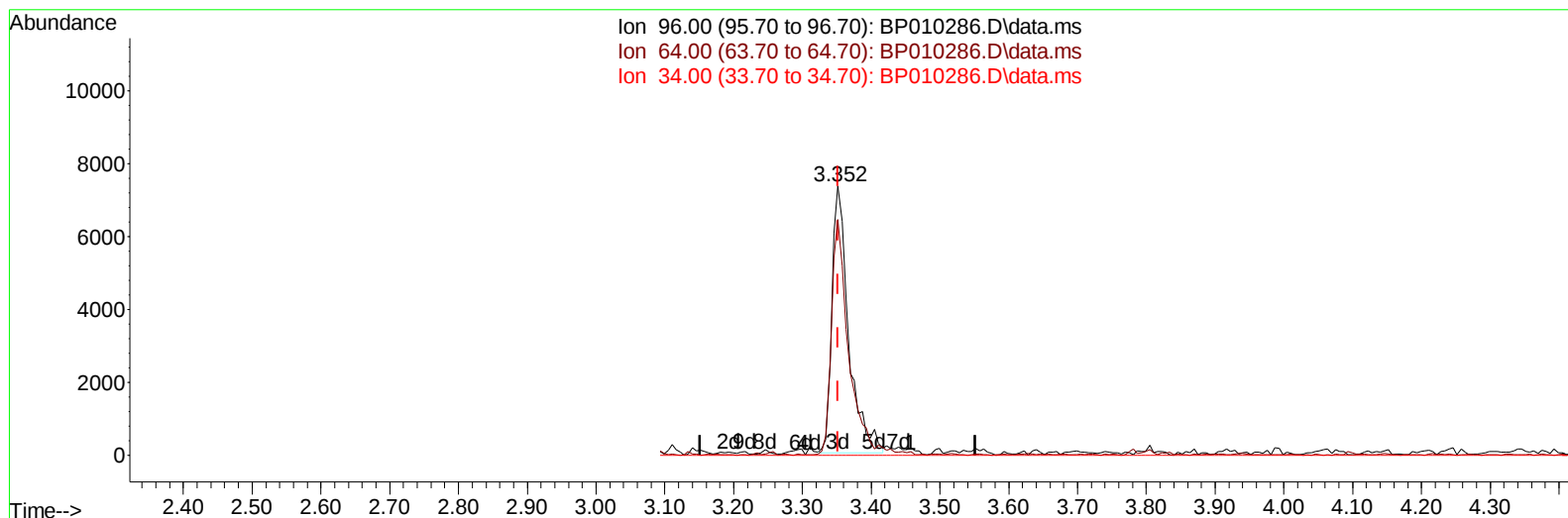
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TIC: BP010286.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.352min (-0.000) 4.34 ng/uL m

response 12393

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	87.59#
34.00	0.00	0.00
0.00	0.00	0.00

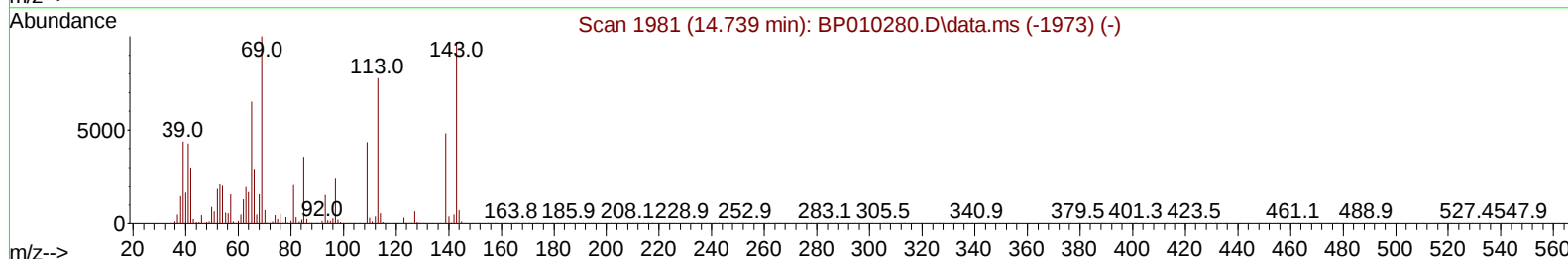
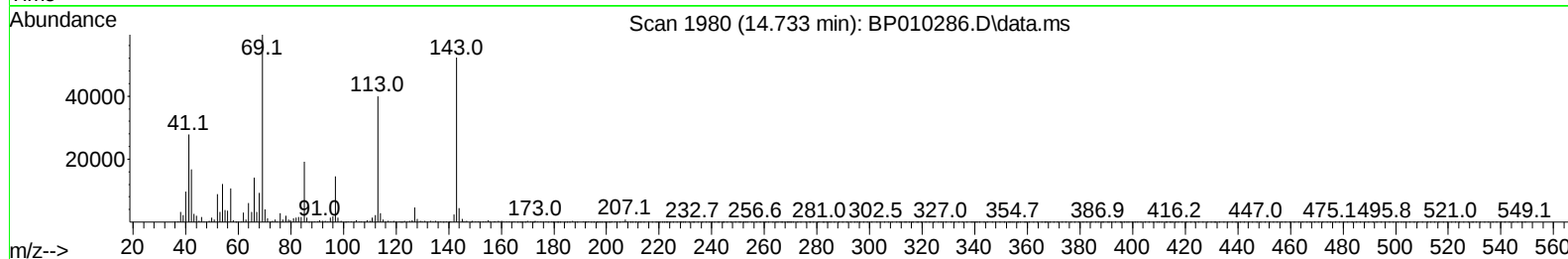
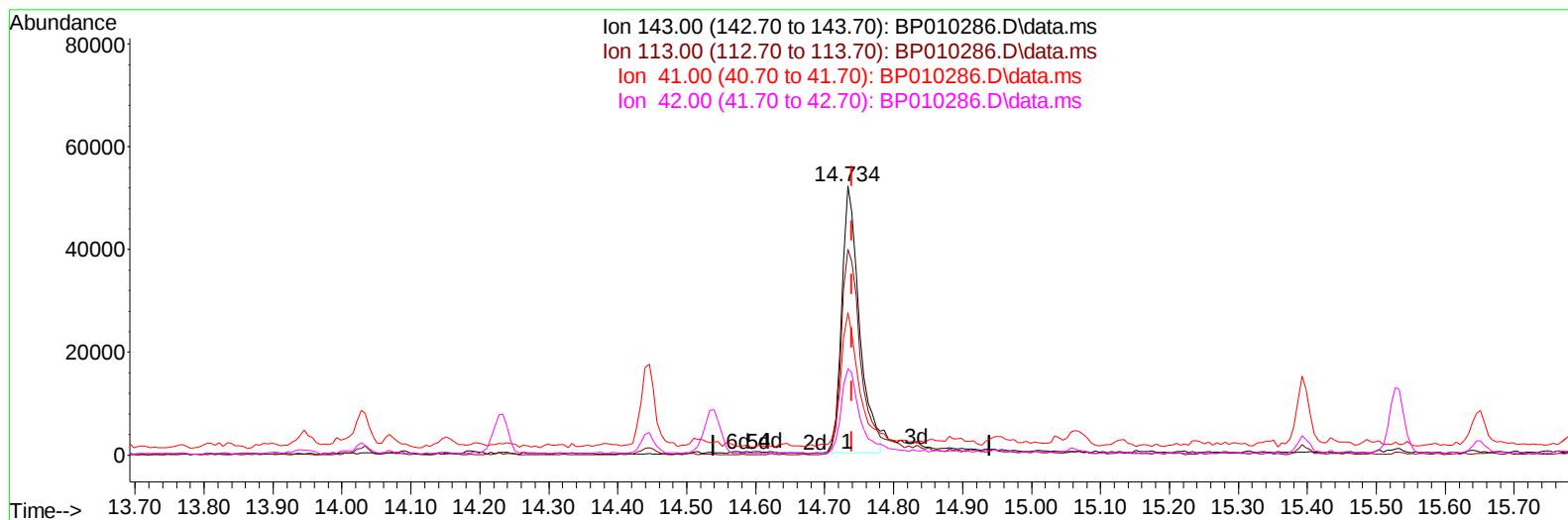
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 Misc :
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Instrument :
 BNA_P
 ClientSampleId :
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Manual IntegrationsAPPROVED

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TIC: BP010286.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.733min (-0.006) 16.12 ng/u1

response 93084

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	76.43#
41.00	23.20	53.03#
42.00	18.00	32.10#

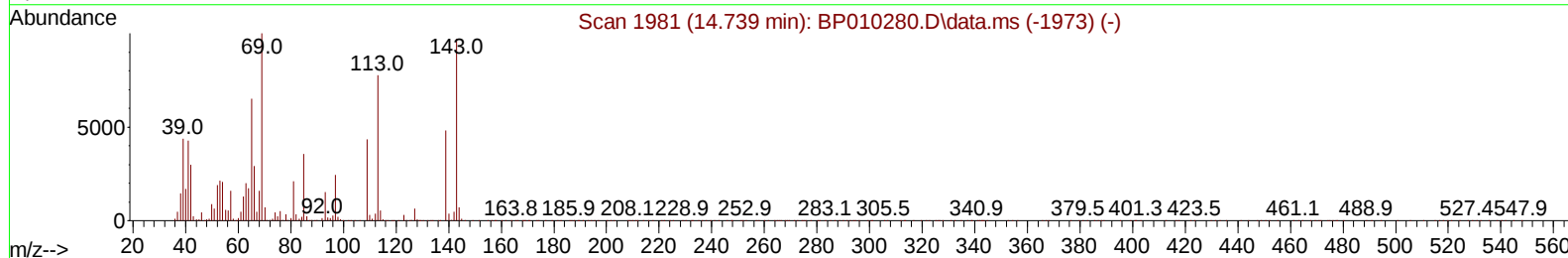
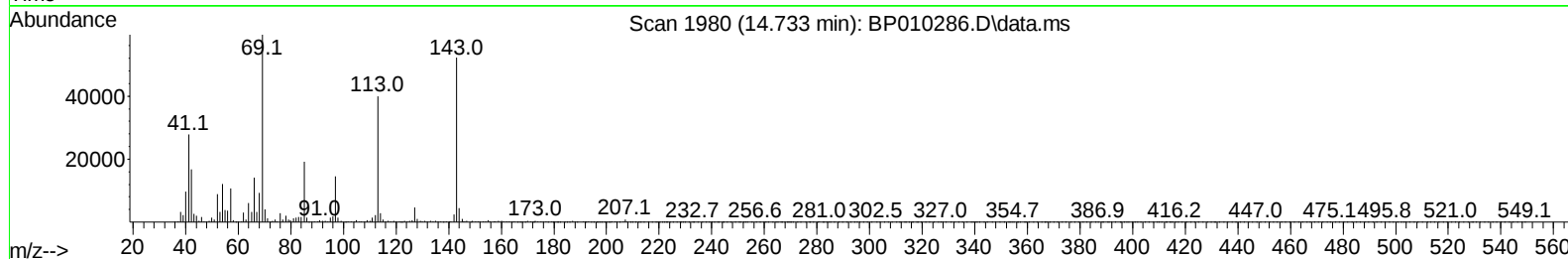
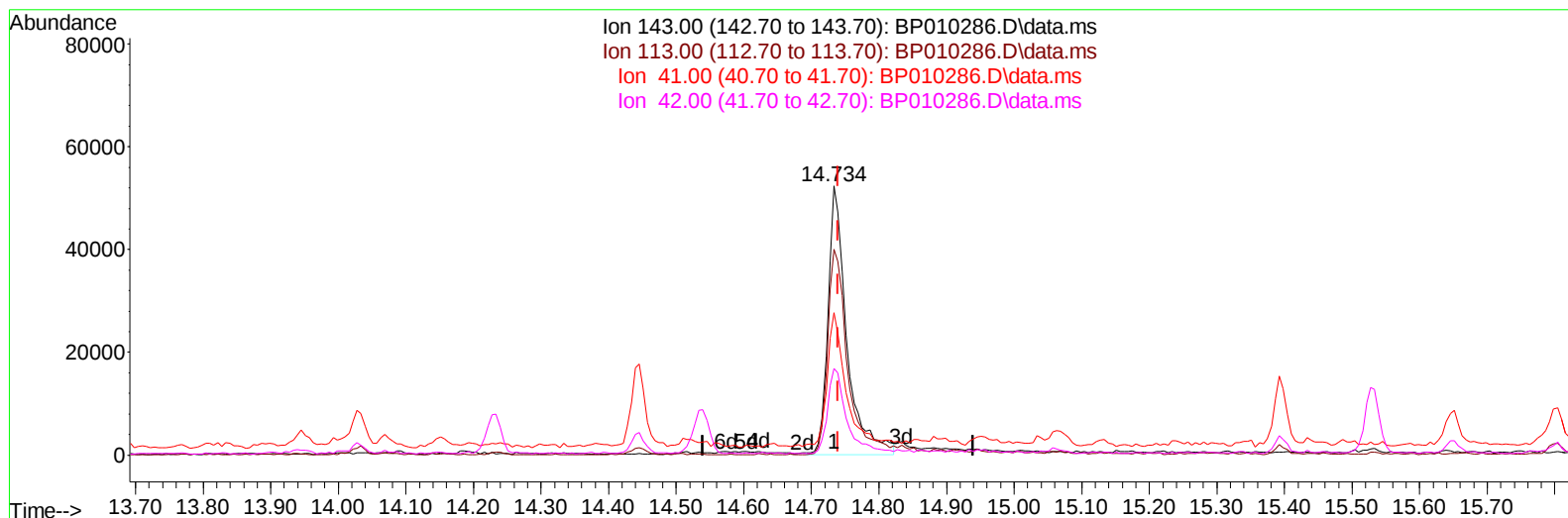
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 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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Manual IntegrationsAPPROVED

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TIC: BP010286.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.733min (-0.006) 17.61 ng/ul m

response 101646

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	76.43#
41.00	23.20	53.03#
42.00	18.00	32.10#

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 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.910	152	121333	20.000	ng/ul	0.00
20) Naphthalene-d8	10.710	136	539789	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.539	164	377779	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.286	188	826403	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.368	240	841205	20.000	ng/ul	# 0.00
88) Perylene-d12	23.804	264	838204	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	12393m	4.339	ng/uL	0.00
4) Pyridine-d5	3.781	84	51668	6.501	ng/ul	0.01
7) Phenol-d5	7.069	99	241563	21.436	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.234	67	173458	24.881	ng/ul	0.00
11) 2-Chlorophenol-d4	7.440	132	197536	23.801	ng/ul	0.00
15) 4-Methylphenol-d8	8.616	113	198424	21.417	ng/ul	0.00
21) Nitrobenzene-d5	9.063	128	105104	24.284	ng/ul	0.00
24) 2-Nitrophenol-d4	9.793	143	112918	22.882	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.334	165	195549	21.772	ng/ul	0.00
31) 4-Chloroaniline-d4	10.845	131	264967	19.983	ng/ul	0.00
46) Dimethylphthalate-d6	13.945	166	747387	24.833	ng/ul	0.00
49) Acenaphthylene-d8	14.234	160	830549	23.146	ng/ul	0.00
54) 4-Nitrophenol-d4	14.733	143	101646m	17.607	ng/ul	0.00
60) Fluorene-d10	15.533	176	654052	25.745	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200	97191	17.174	ng/ul	0.00
73) Anthracene-d10	17.386	188	1079328	26.670	ng/ul	0.00
81) Pyrene-d10	19.621	212	1428043	30.720	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.645	264	1316652	29.454	ng/ul	-0.01

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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