

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041724\
Data File : BM045332.D
Acq On : 18 Apr 2024 00:59
Operator : MA/JU
Sample : P2145-03
Misc :
ALS Vial : 24 Sample Multiplier: 1

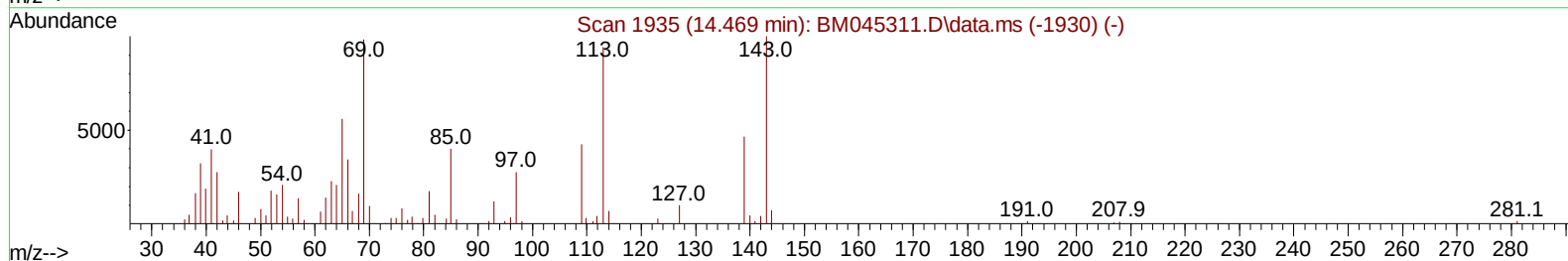
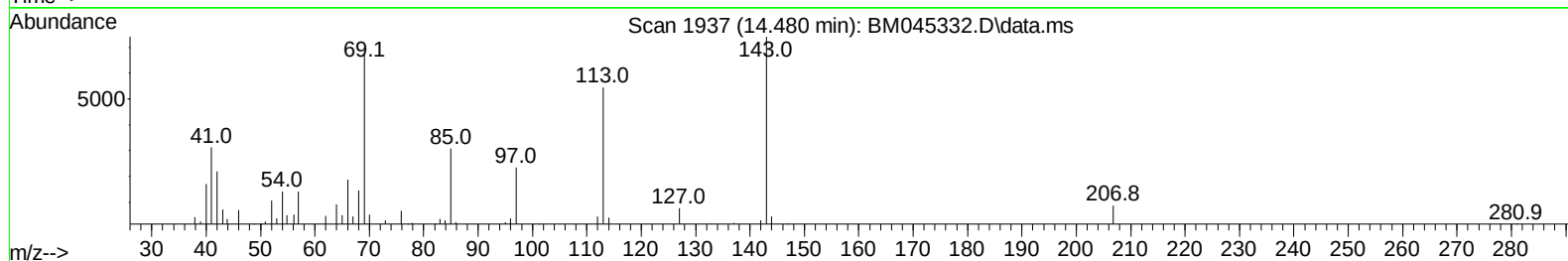
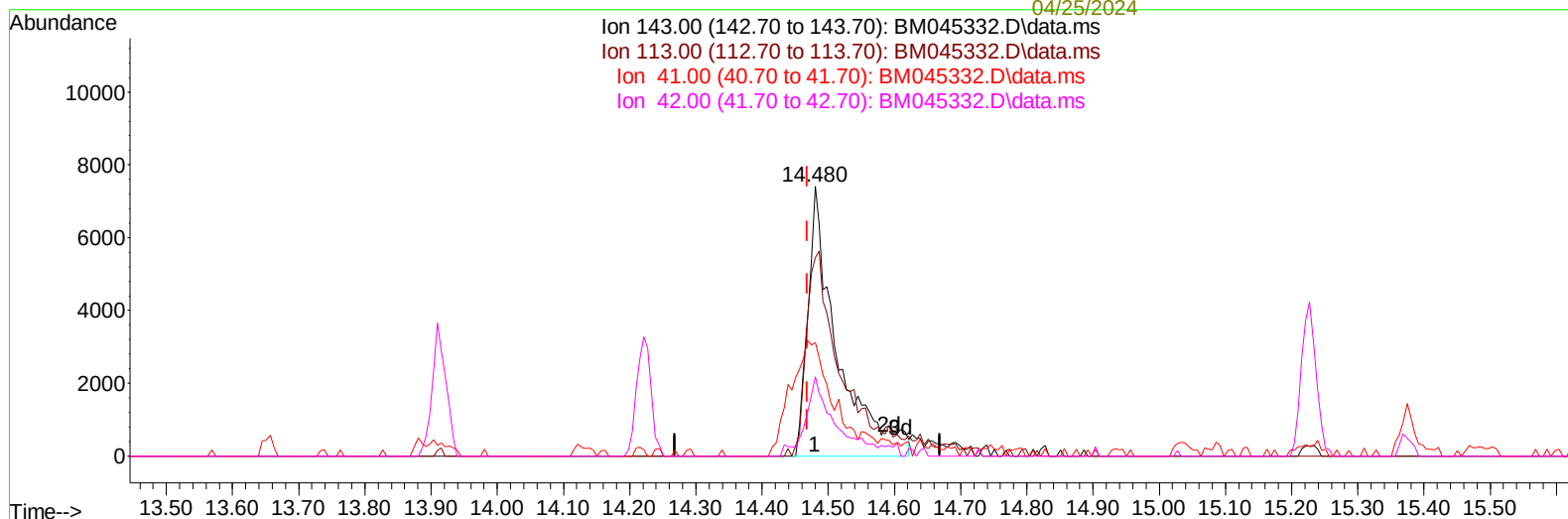
Instrument :
BNA_M
ClientSampleId :
E07X1

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 04/19/2024
Supervised By :mohammad ahmed 04/25/2024

Supervised By :mohammad
ahmed
04/25/2024

Quant Time: Apr 18 02:35:18 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Apr 17 12:42:01 2024
Response via : Initial Calibration



TIC: BM045332.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.480min (+ 0.012) 17.30 ng/ul m

response 22993

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.80	73.58
41.00	36.50	42.13
42.00	23.30	29.46#

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Quant Time: Apr 18 02:36:26 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.569	152	39971	20.000	ng/ul	0.00
20) Naphthalene-d8	10.339	136	153857	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.221	164	91480	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.986	188	180421	20.000	ng/ul	0.00
79) Chrysene-d12	21.197	240	185520	20.000	ng/ul	0.00
88) Perylene-d12	23.397	264	245861	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	4309	4.026	ng/uL	0.00
4) Pyridine-d5	3.575	84	51304	17.625	ng/ul	0.00
7) Phenol-d5	6.757	99	76899	22.694	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.928	67	49905	24.719	ng/ul	0.00
11) 2-Chlorophenol-d4	7.104	132	67238	25.327	ng/ul	0.00
15) 4-Methylphenol-d8	8.281	113	66464	25.058	ng/ul	0.00
21) Nitrobenzene-d5	8.734	128	31597	26.736	ng/ul	0.00
24) 2-Nitrophenol-d4	9.445	143	33952	27.218	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.975	165	61195	26.757	ng/ul	0.00
31) 4-Chloroaniline-d4	10.504	131	86295	23.699	ng/ul	0.00
46) Dimethylphthalate-d6	13.651	166	191588	30.079	ng/ul	0.00
49) Acenaphthylene-d8	13.916	160	223664	29.724	ng/ul	0.00
54) 4-Nitrophenol-d4	14.480	143	22993m	17.297	ng/ul	0.01
60) Fluorene-d10	15.227	176	167528	29.502	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.374	200	18607	17.394	ng/ul	0.00
73) Anthracene-d10	17.086	188	252005	31.049	ng/ul	0.00
81) Pyrene-d10	19.392	212	295897	32.392	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.256	264	376874	31.422	ng/ul	0.00

Target Compounds Qvalue

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

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