

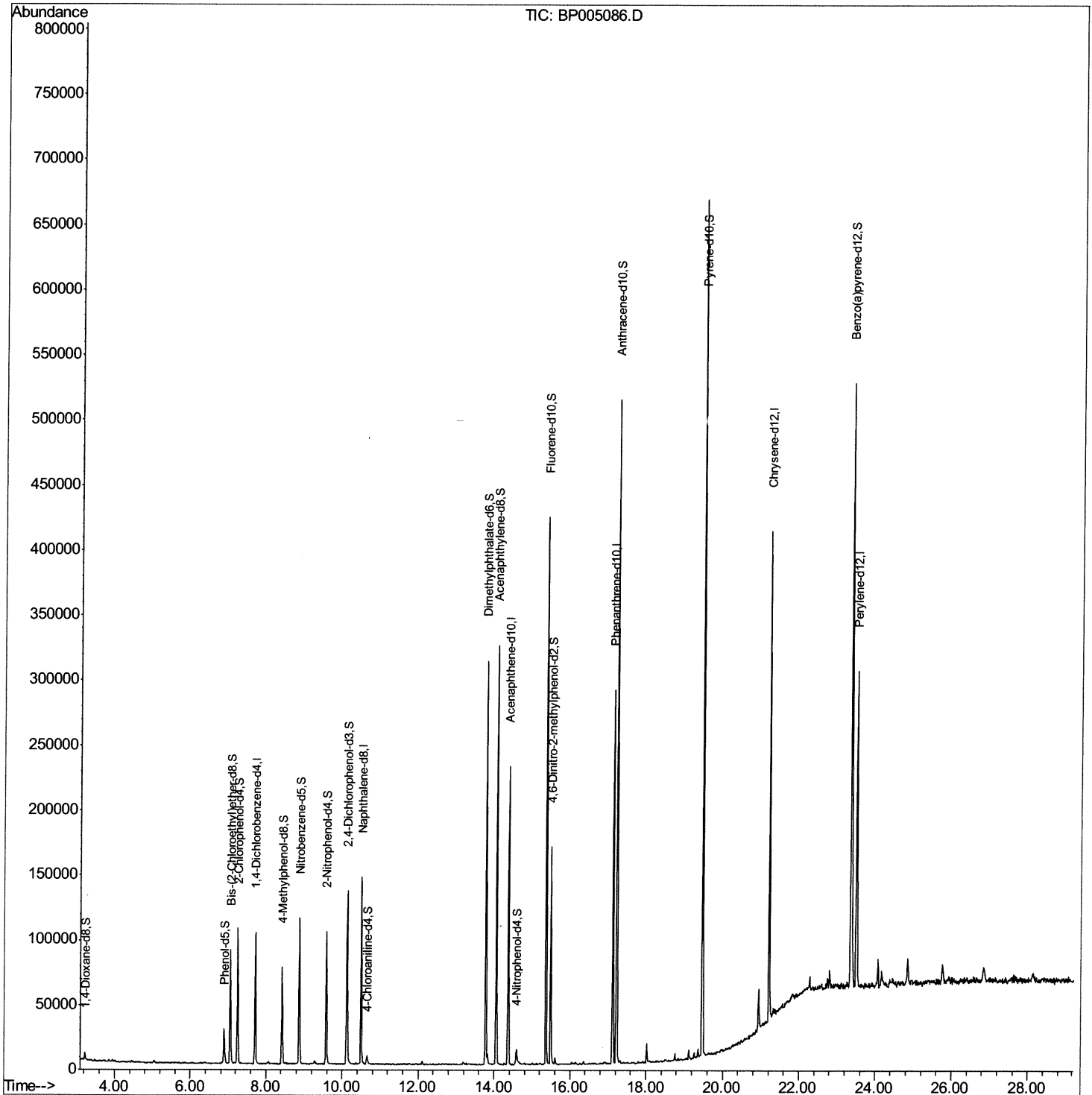
Data Path : Z:\svoasrv\HPCHEM1\BNA P\Data\BP032921\
Data File : BP005086.D
Acq On : 29 Mar 2021 21:54
Operator : CG/JU
Sample : M1783-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0G08

Manual Integrations
APPROVED

mohammad
3/30/2021 4:12:28 PM

Quant Time: Mar 30 02:04:18 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP032921MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Mar 30 00:27:26 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

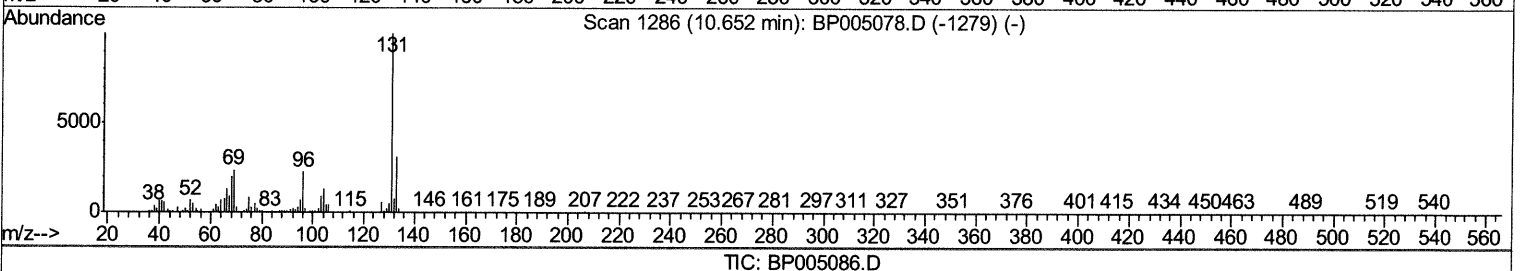
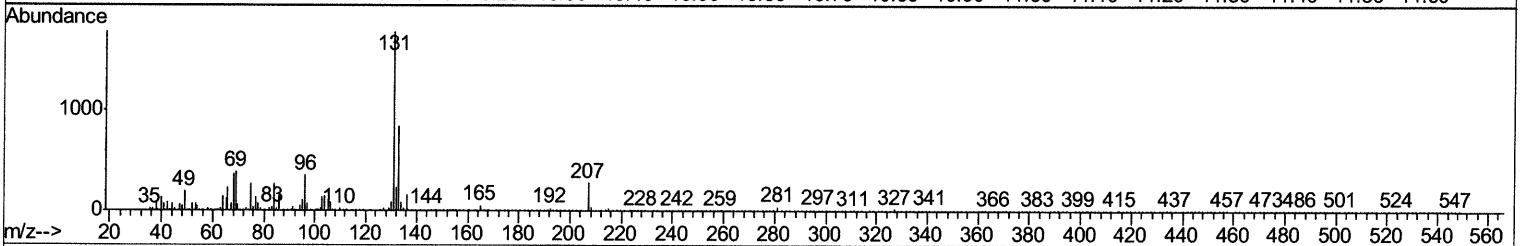
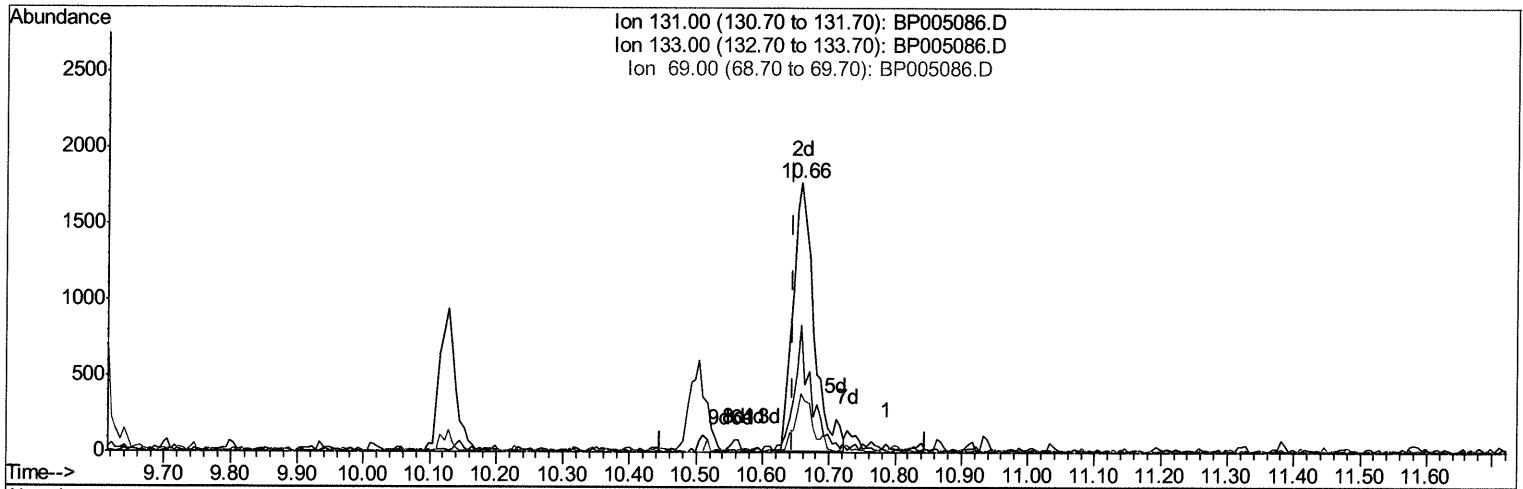
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(29) 4-Chloroaniline-d4 (S)

10.657min (+0.012) 2.01ng/ul m 04/01/2021

response 3888

Ion	Exp%	Act%
131.00	100	100
133.00	31.40	47.23#
69.00	21.50	21.90
0.00	0.00	0.00

Quantitation Report (Qedit)

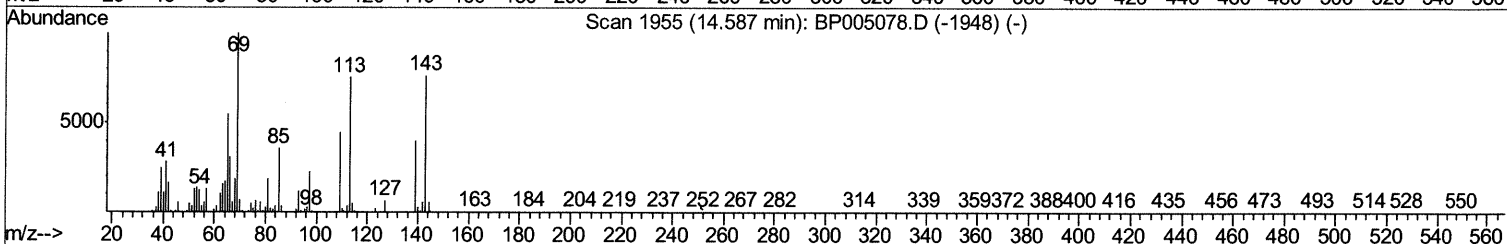
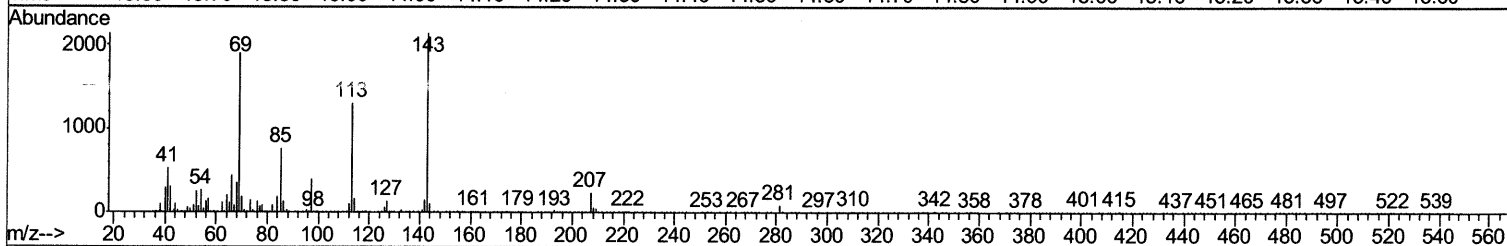
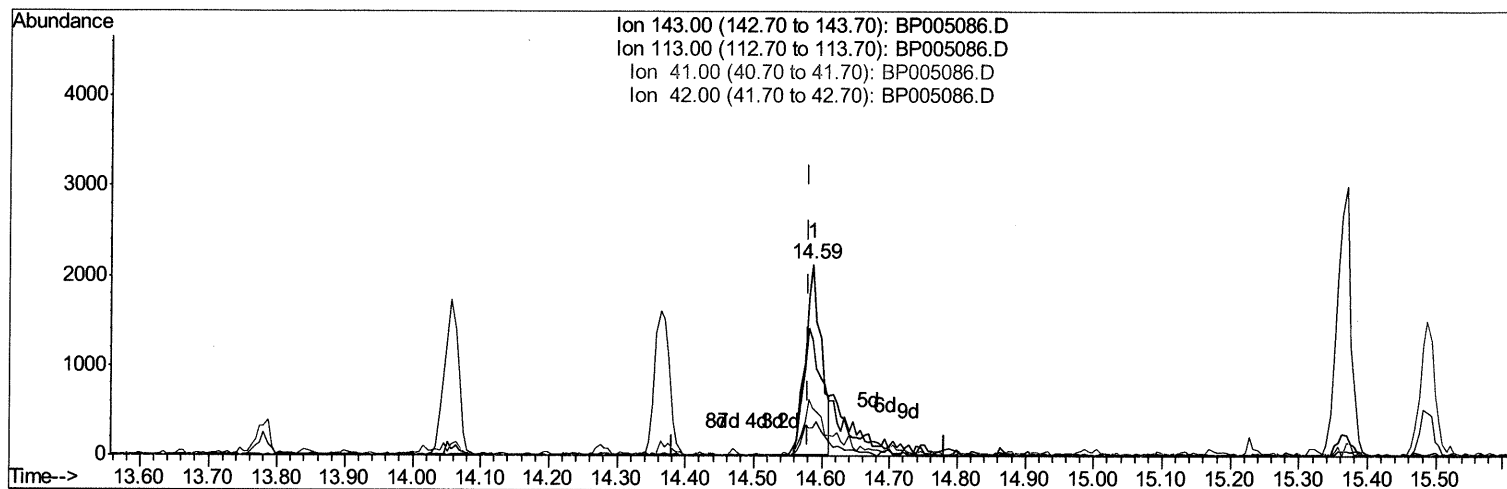
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TIC: BP005086.D

(52) 4-Nitrophenol-d4 (S)

14.587min (+0.006) 3.96ng/ul

response 3503

Ion	Exp%	Act%
143.00	100	100
113.00	81.90	61.15#
41.00	27.10	24.96
42.00	18.80	14.45#

Quantitation Report (Qedit)

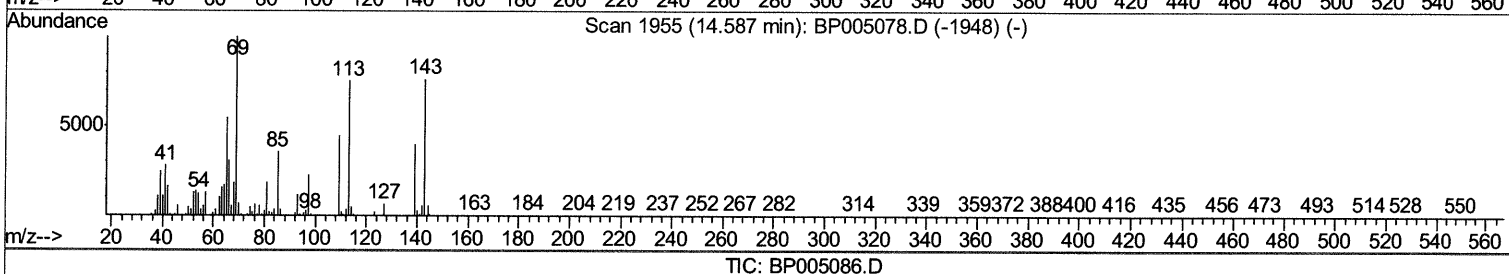
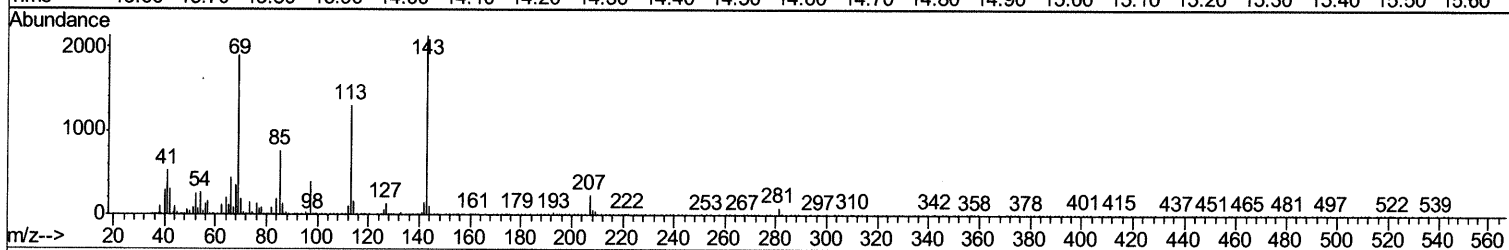
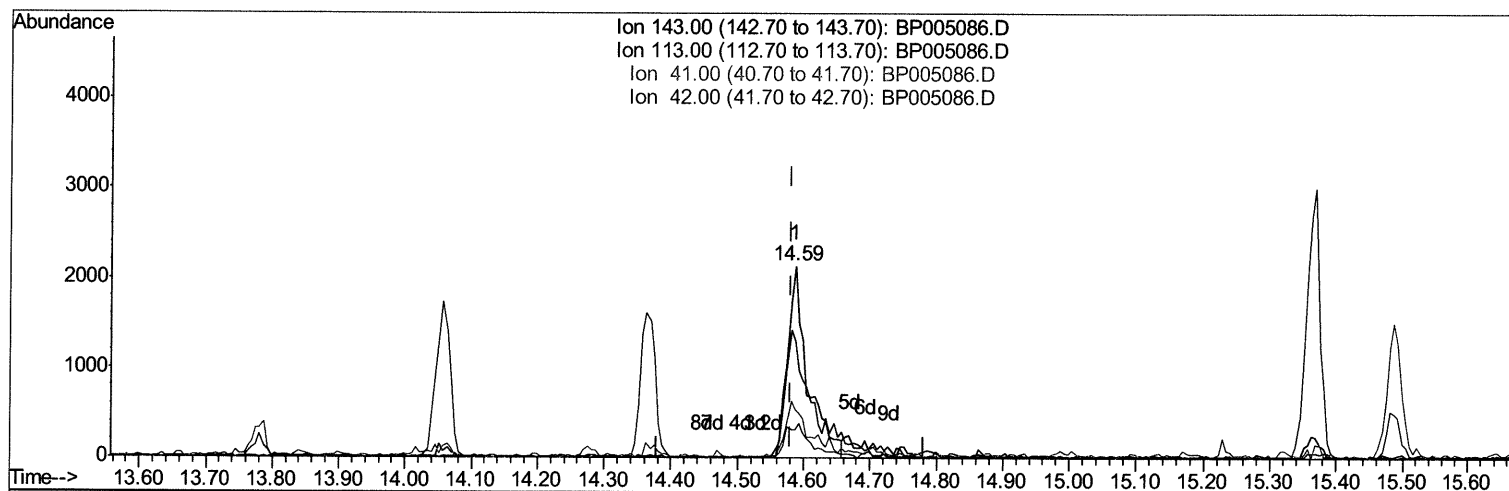
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Operator : CG/JU
Sample : M1783-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0G08

Manual Integrations
APPROVED

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(52) 4-Nitrophenol-d4 (S)

14.587min (+0.006) 5.13ng/ul m

04/01/21 JU

response 4540

Ion	Exp%	Act%
143.00	100	100
113.00	81.90	61.15#
41.00	27.10	24.96
42.00	18.80	14.45#

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	26206	20.00	ng/ul	0.00
18) Naphthalene-d8	10.50	136	110005	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.37	164	74558	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.13	188	166139	20.00	ng/ul	0.00
78) Chrysene-d12	21.22	240	172867	20.00	ng/ul	0.00
86) Perylene-d12	23.50	264	168410	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.22	96	3251	4.69	ng/uL	0.00
5) Phenol-d5	6.89	99	15027	6.66	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.05	67	38216	32.34	ng/ul	0.00
9) 2-Chlorophenol-d4	7.25	132	41104	25.85	ng/ul	0.00
13) 4-Methylphenol-d8	8.42	113	27259	15.81	ng/ul	0.00
19) Nitrobenzene-d5	8.87	128	25708	31.66	ng/ul	0.00
22) 2-Nitrophenol-d4	9.59	143	28558	31.80	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.13	165	46452	26.51	ng/ul	0.00
29) 4-Chloroaniline-d4	10.66	131	3888m>	2.01	ng/ul >	0.01 04/01/21JU
44) Dimethylphthalate-d6	13.78	166	182899	33.78	ng/ul	0.00
47) Acenaphthylene-d8	14.06	160	215662	32.14	ng/ul	0.00
52) 4-Nitrophenol-d4	14.59	143	4540m>	5.13	ng/ul >	0.00 04/01/21JU
58) Fluorene-d10	15.36	176	160310	34.25	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.49	200	35552	32.40	ng/ul	0.00
71) Anthracene-d10	17.23	188	287060	38.38	ng/ul	0.00
79) Pyrene-d10	19.47	212	324909	39.28	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.36	264	343580	39.72	ng/ul	0.00

Target Compounds Ovalue

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