

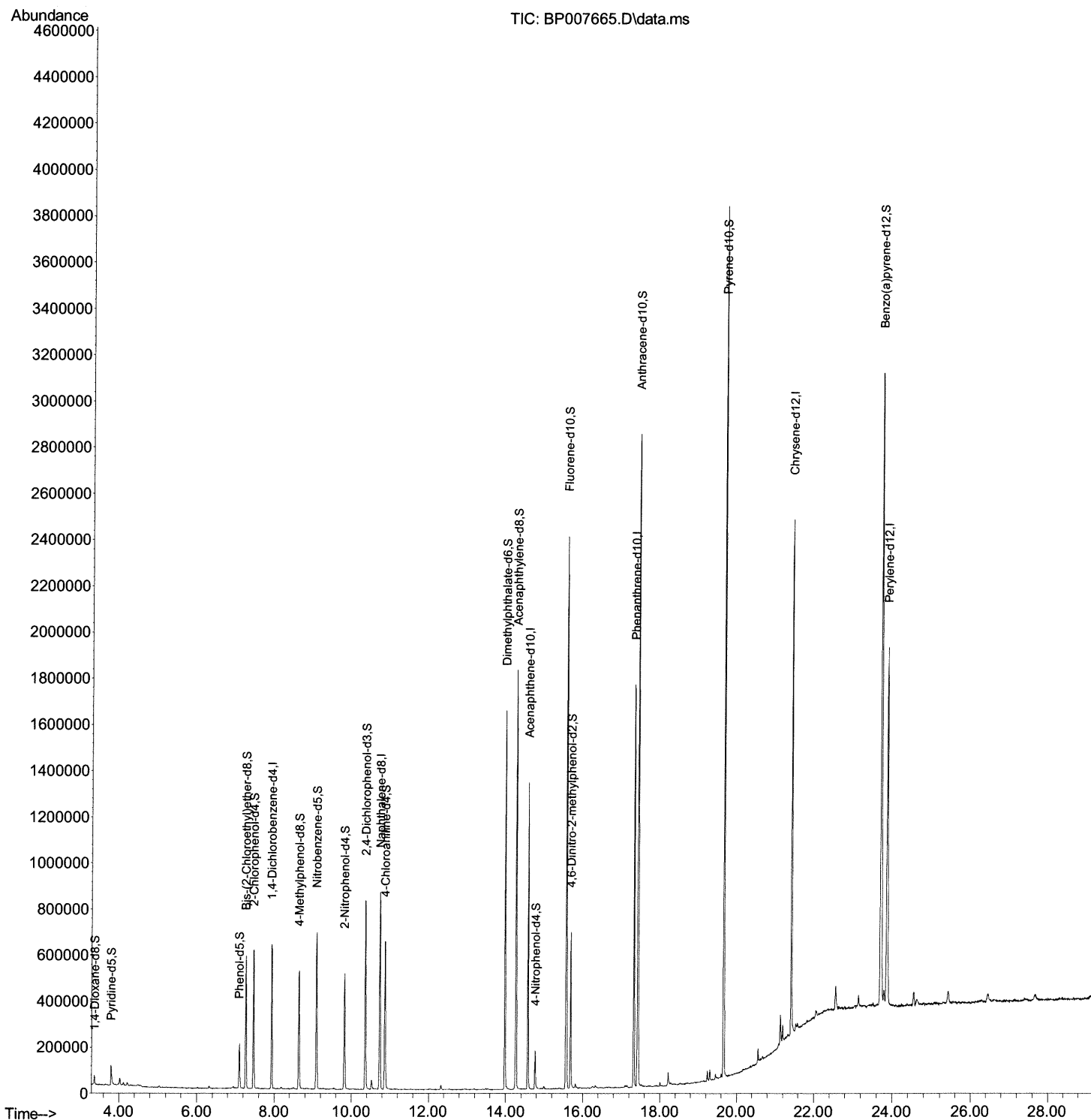
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
Data File : BP007665.D
Acq On : 26 Oct 2021 05:00
Operator : CG/JU
Sample : M4291-10
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
P022-TW001-02

Manual Integrations
APPROVED

mohammad
10/26/2021 11:35:00 AM

Quant Time: Oct 26 05:49:34 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Oct 25 16:57:16 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

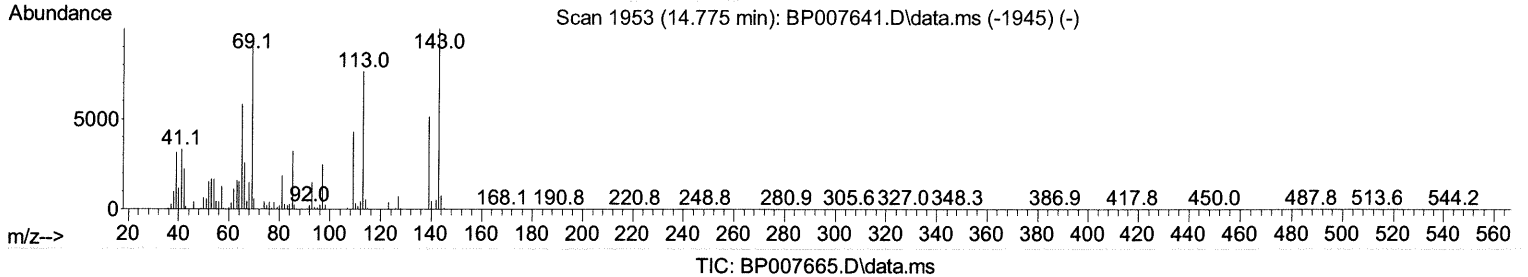
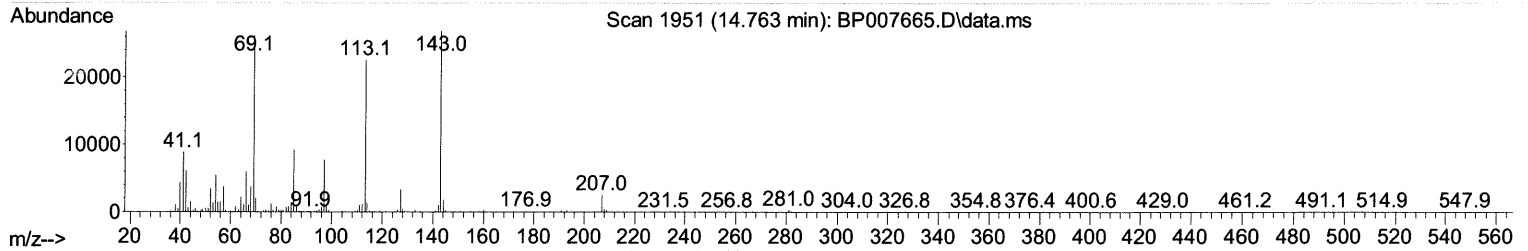
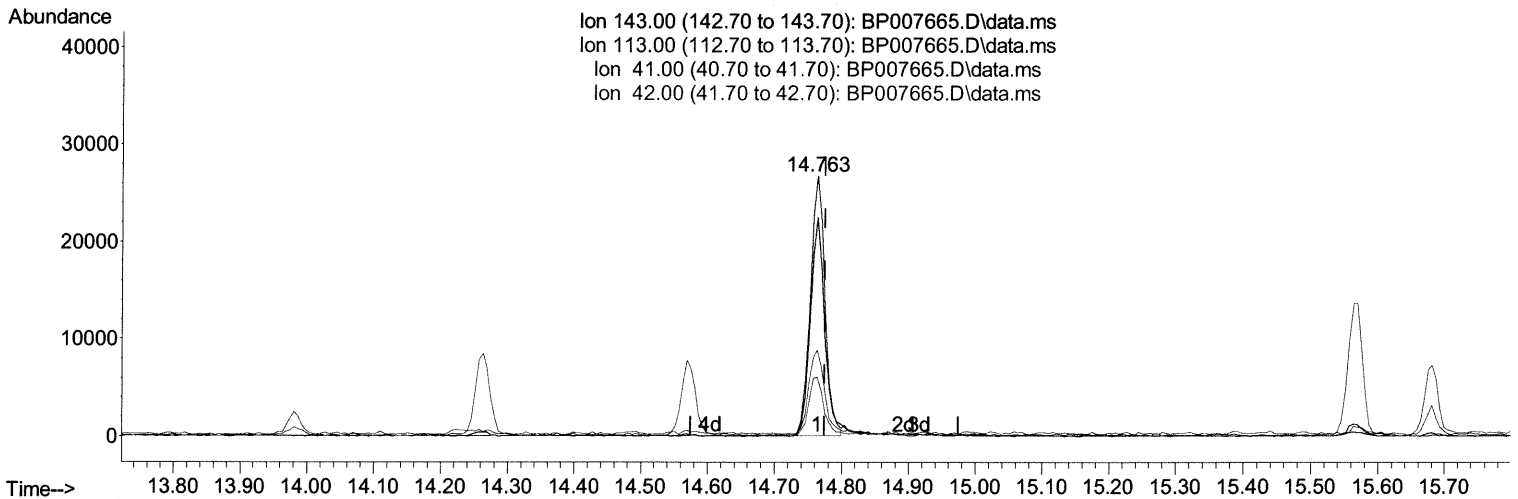
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(54) 4-Nitrophenol-d4 (S)

14.763min (-0.012) 7.01 ng/ul

response 40108

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	79.90	84.14
41.00	39.60	33.06
42.00	25.40	22.79

Quantitation Report (Qedit)

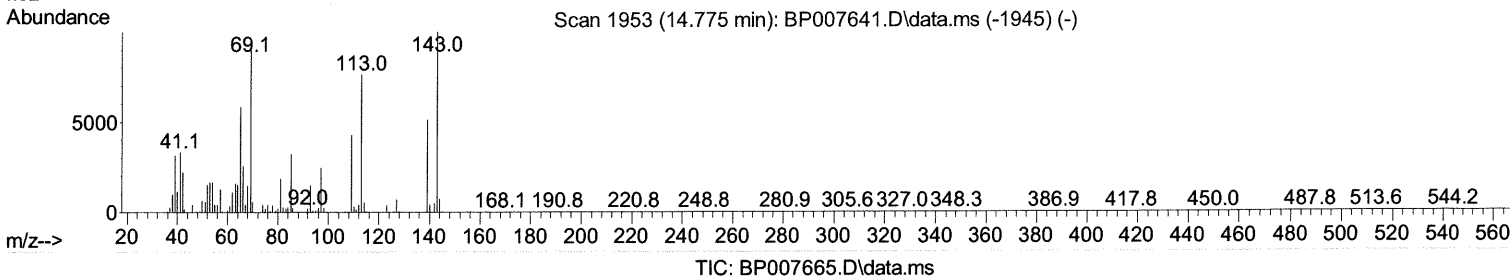
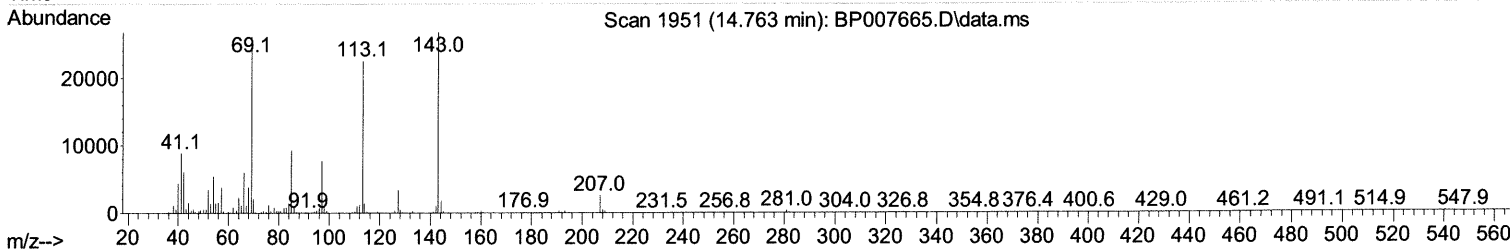
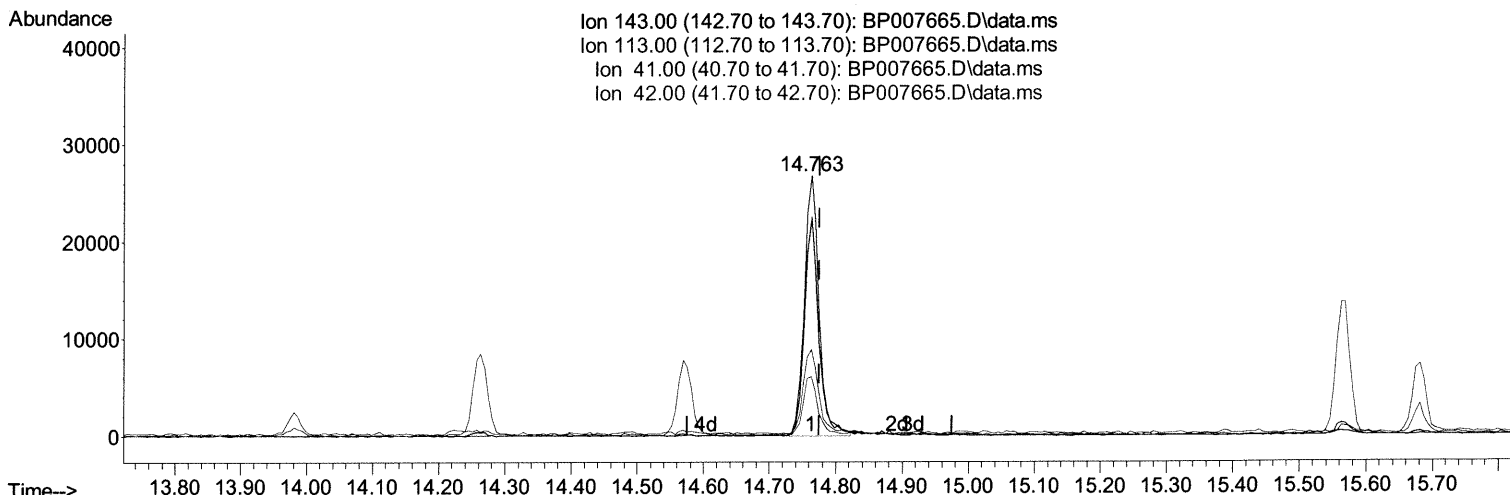
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(54) 4-Nitrophenol-d4 (S)

14.763min (-0.012) 7.18 ng/ul m 10/27/21ju

response 41078

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	79.90	84.14
41.00	39.60	33.06
42.00	25.40	22.79

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	172697	20.000	ng/ul	0.00
20) Naphthalene-d8	10.740	136	694229	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.575	164	460866	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.328	188	1026975	20.000	ng/ul	0.00
79) Chrysene-d12	21.416	240	1133911	20.000	ng/ul	0.00
88) Perylene-d12	23.857	264	1183430	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	21166	5.567	ng/uL	0.00
4) Pyridine-d5	3.793	84	52051	4.527	ng/ul	0.00
7) Phenol-d5	7.099	99	109561	8.352	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.264	67	249749	34.293	ng/ul	0.00
11) 2-Chlorophenol-d4	7.464	132	284311	27.868	ng/ul	-0.01
15) 4-Methylphenol-d8	8.646	113	191066	18.602	ng/ul	0.00
21) Nitrobenzene-d5	9.093	128	159934	33.806	ng/ul	0.00
24) 2-Nitrophenol-d4	9.816	143	164052	31.749	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.358	165	308177	29.301	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.875	131	348087	25.030	ng/ul	0.00
46) Dimethylphthalate-d6	13.981	166	1113656	35.674	ng/ul	-0.01
49) Acenaphthylene-d8	14.263	160	1276782	33.369	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	41078m >	7.178	ng/ul >	-0.01
60) Fluorene-d10	15.569	176	964799	36.660	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	151528	27.584	ng/ul	0.00
73) Anthracene-d10	17.428	188	1663453	39.414	ng/ul	0.00
81) Pyrene-d10	19.657	212	2062563	39.070	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.698	264	2139501	36.802	ng/ul	-0.01

Target Compounds Qvalue

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