

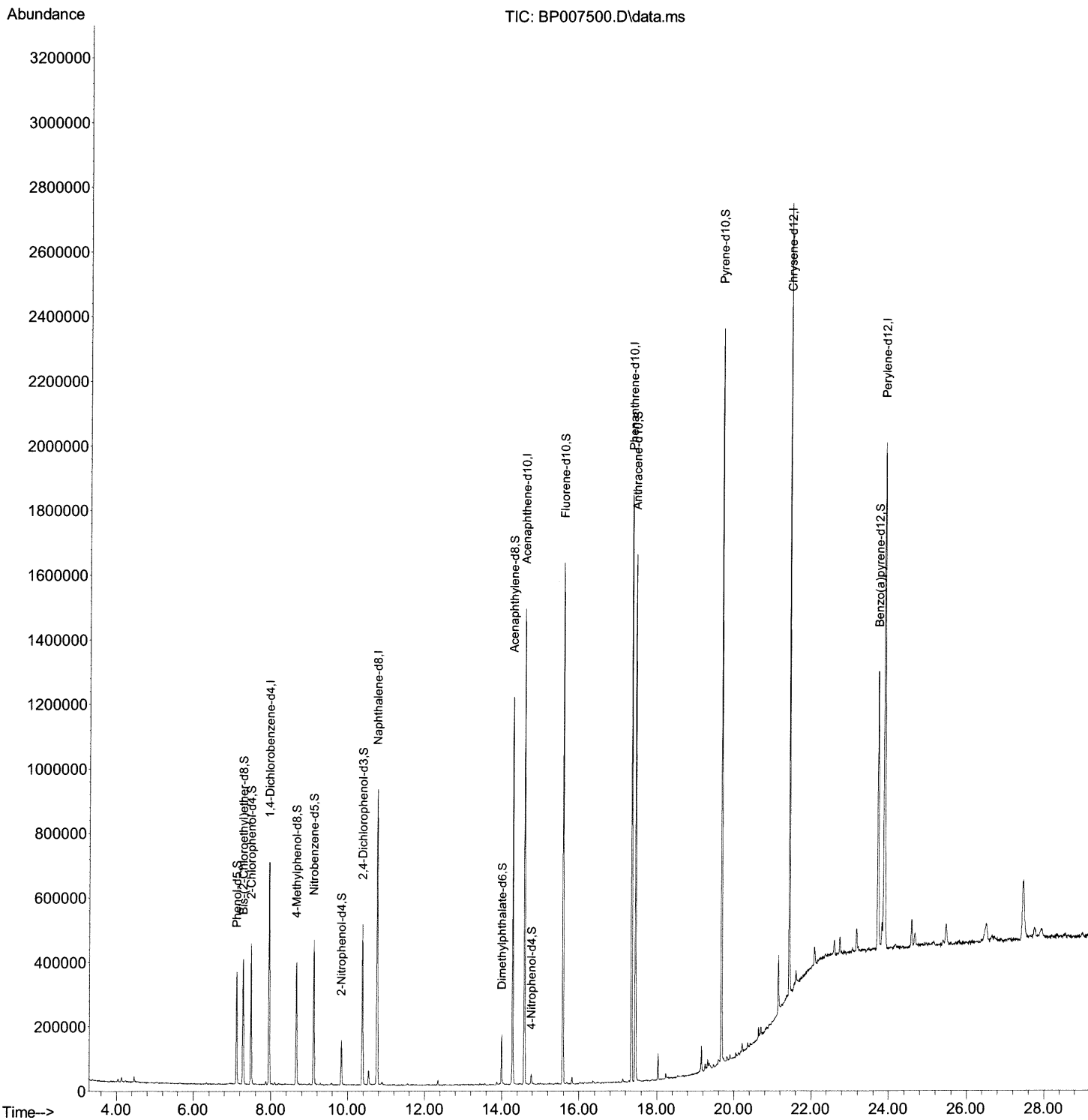
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007500.D
Acq On : 19 Oct 2021 18:46
Operator : CG/JU
Sample : M4196-01
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JDGA0

Manual Integrations
APPROVED

mohammad
10/20/2021 1:56:43 PM

Quant Time: Oct 20 01:04:54 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Oct 18 11:26:58 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

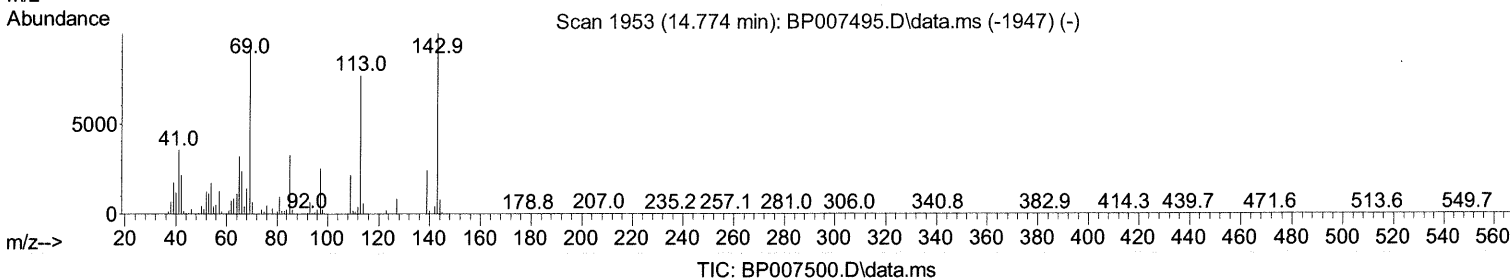
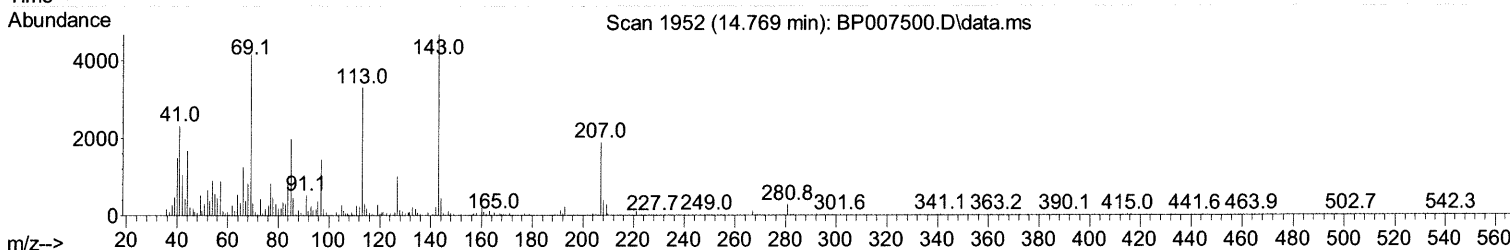
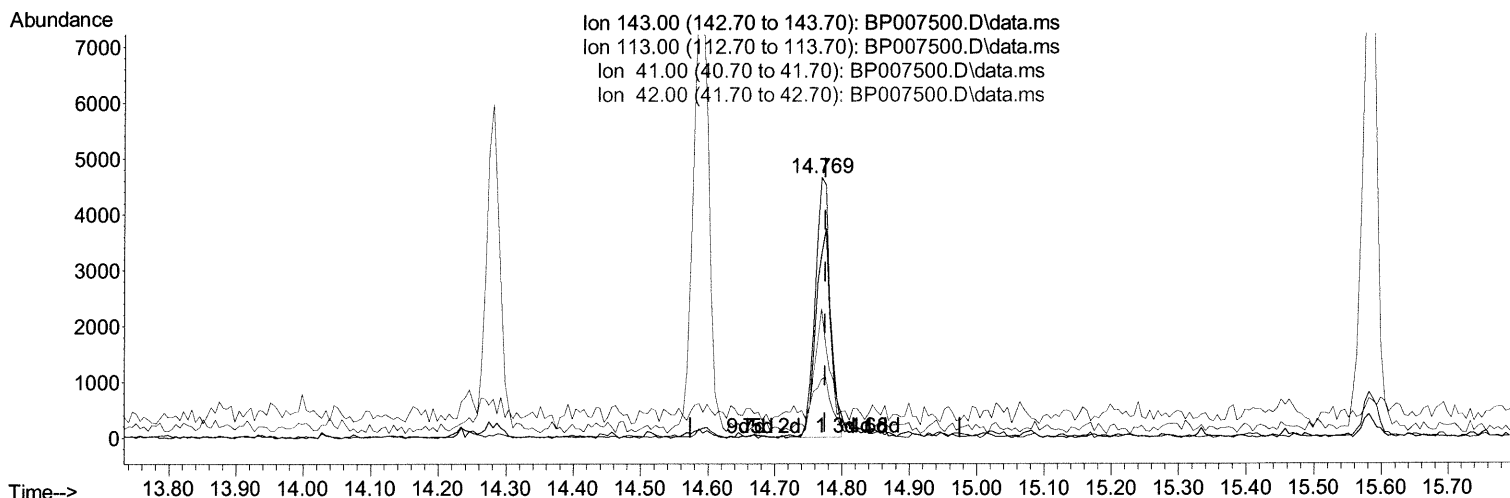
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(54) 4-Nitrophenol-d4 (S)

14.769min (-0.006) 1.26 ng/ul

response 7249

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	79.90	70.71
41.00	39.60	49.66#
42.00	25.40	22.23

Quantitation Report (Qedit)

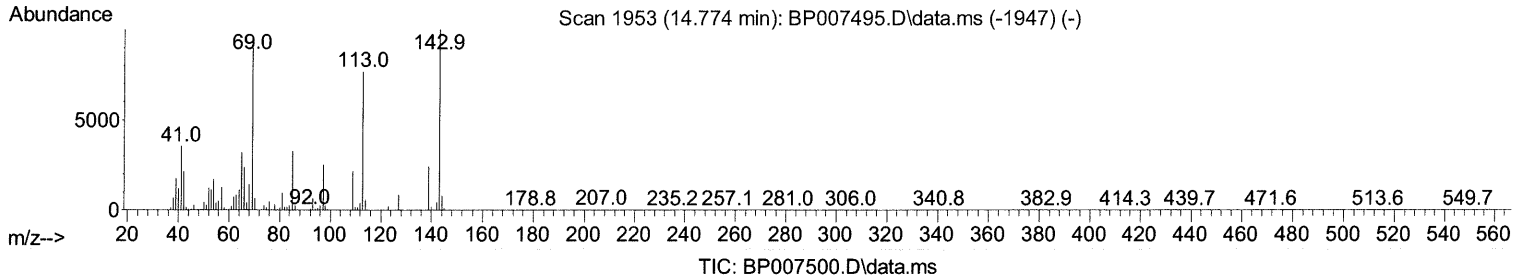
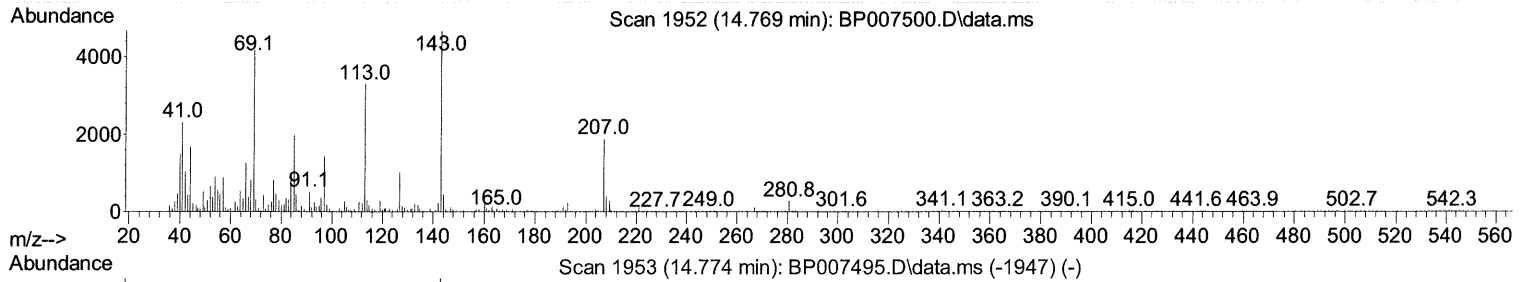
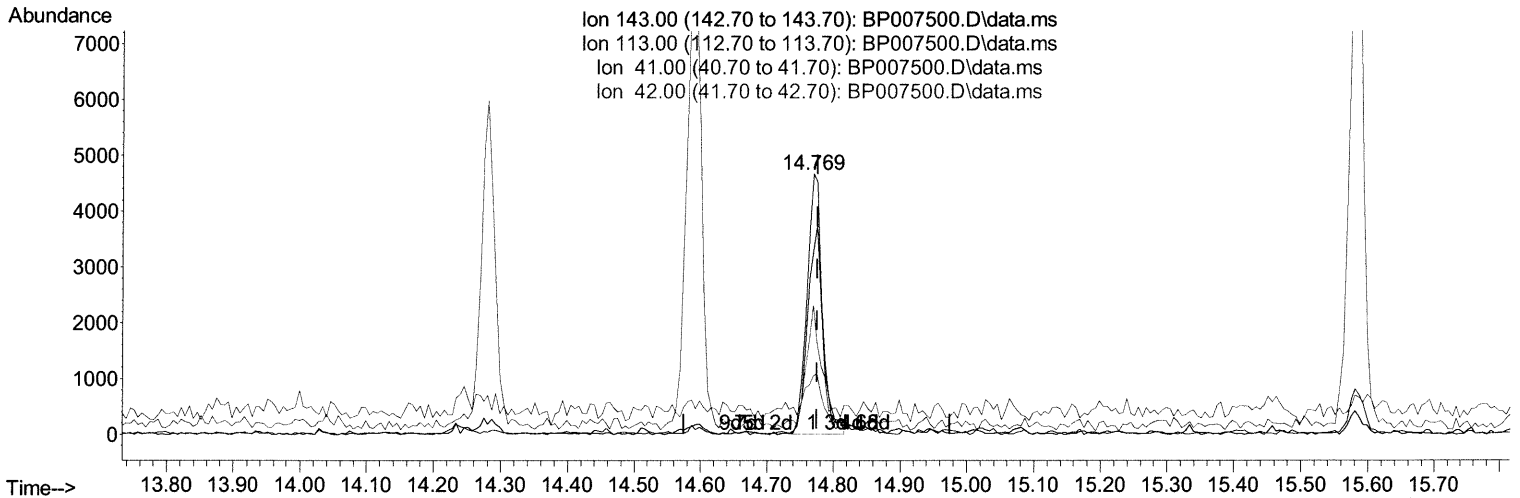
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(54) 4-Nitrophenol-d4 (S)

14.769min (-0.006) 1.29 ng/ul m 10/21/21 JU

response 7460

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	79.90	70.71
41.00	39.60	49.66#
42.00	25.40	22.23

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
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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	7.958	152	194973	20.000	ng/ul	0.00
20) Naphthalene-d8	10.757	136	790066	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.587	164	519825	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.339	188	1132776	20.000	ng/ul	-0.01
79) Chrysene-d12	21.427	240	1190241	20.000	ng/ul	-0.01
88) Perylene-d12	23.880	264	1224451	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	2993	0.606	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.111	99	200706	13.181	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	169137	18.624	ng/ul	0.00
11) 2-Chlorophenol-d4	7.481	132	202966	16.944	ng/ul	0.00
15) 4-Methylphenol-d8	8.657	113	141510	12.065	ng/ul	0.00
21) Nitrobenzene-d5	9.110	128	108292	21.659	ng/ul	0.00
24) 2-Nitrophenol-d4	9.834	143	44455	9.519	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	181391	15.924	ng/ul	0.00
31) 4-Chloroaniline-d4	10.887	131	3417	0.222	ng/ul	0.00
46) Dimethylphthalate-d6	13.998	166	104051	2.893	ng/ul	0.00
49) Acenaphthylene-d8	14.281	160	854384	18.904	ng/ul	0.00
54) 4-Nitrophenol-d4	14.769	143	7460m>	1.293	ng/ul>	0.00 10/21/21 JU
60) Fluorene-d10	15.581	176	633510	21.049	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.692	200	1137	0.250	ng/ul	-0.01
73) Anthracene-d10	17.439	188	964154	20.042	ng/ul	-0.01
81) Pyrene-d10	19.675	212	1180774	19.844	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.721	264	677138	11.114	ng/ul	-0.01

Target Compounds Qvalue

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