

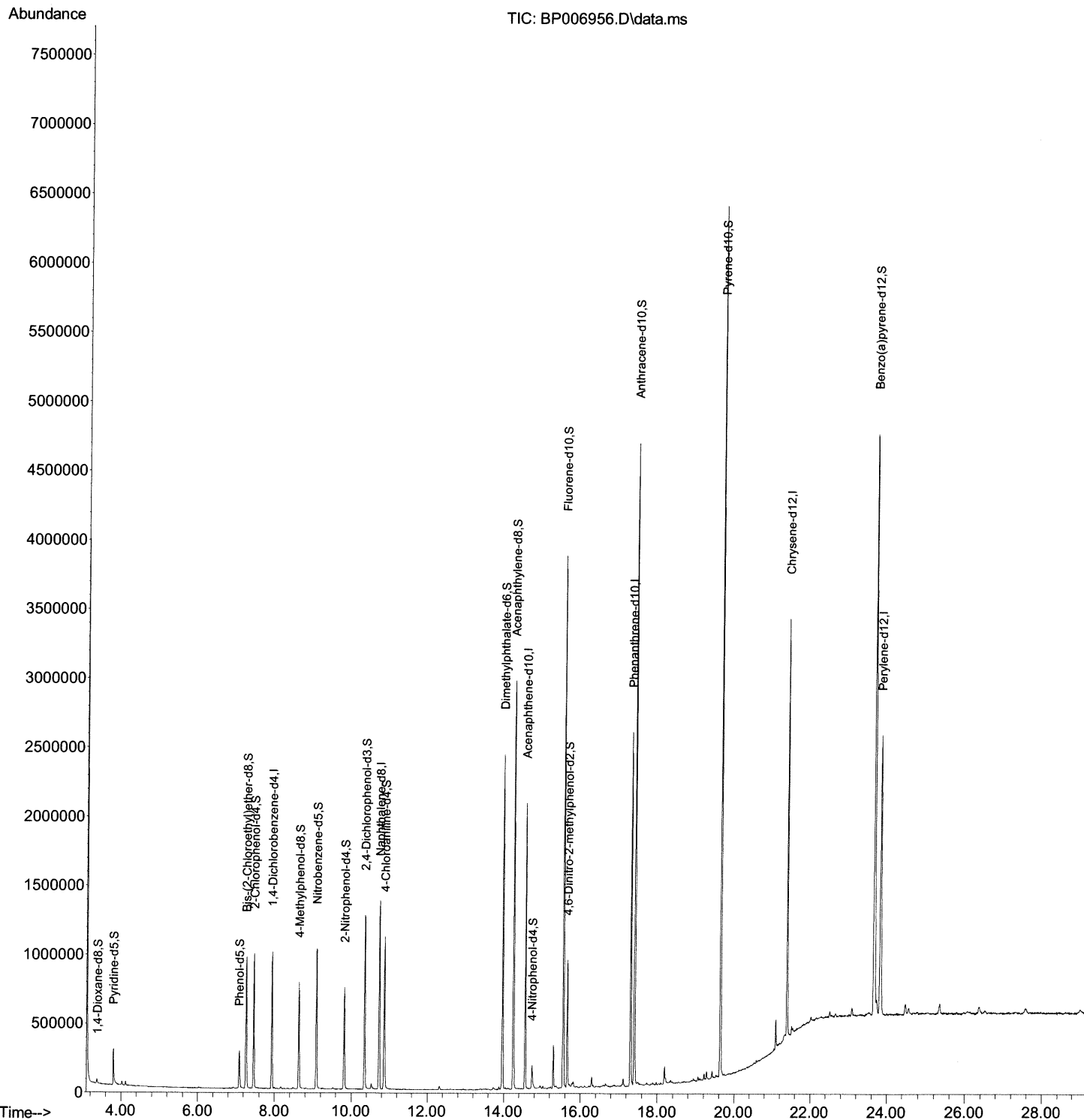
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090921\
Data File : BP006956.D
Acq On : 10 Sep 2021 23:44
Operator : CG/JU
Sample : M3651-12
Misc :
ALS Vial : 56 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBKM6

Manual Integrations
APPROVED

mohammad
9/13/2021 10:22:31 AM

Quant Time: Sep 11 00:35:47 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP090821.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Sep 09 11:07:24 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

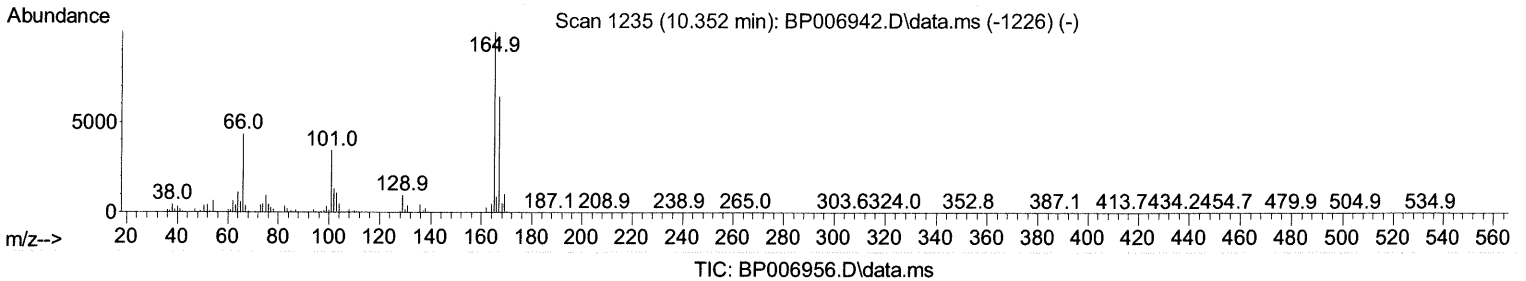
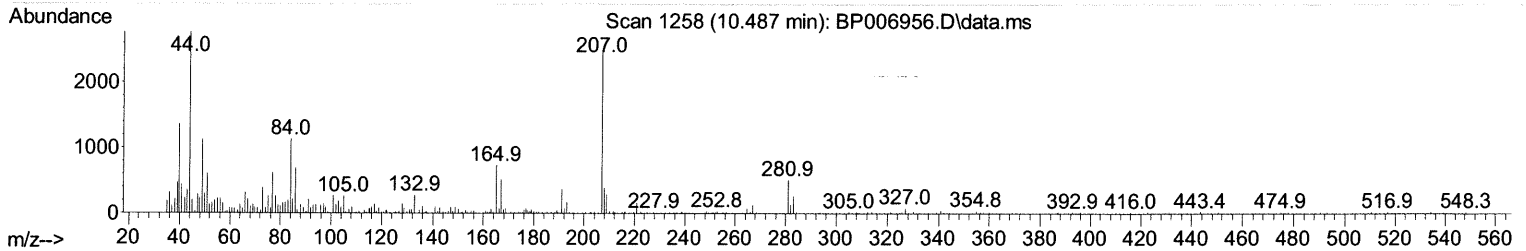
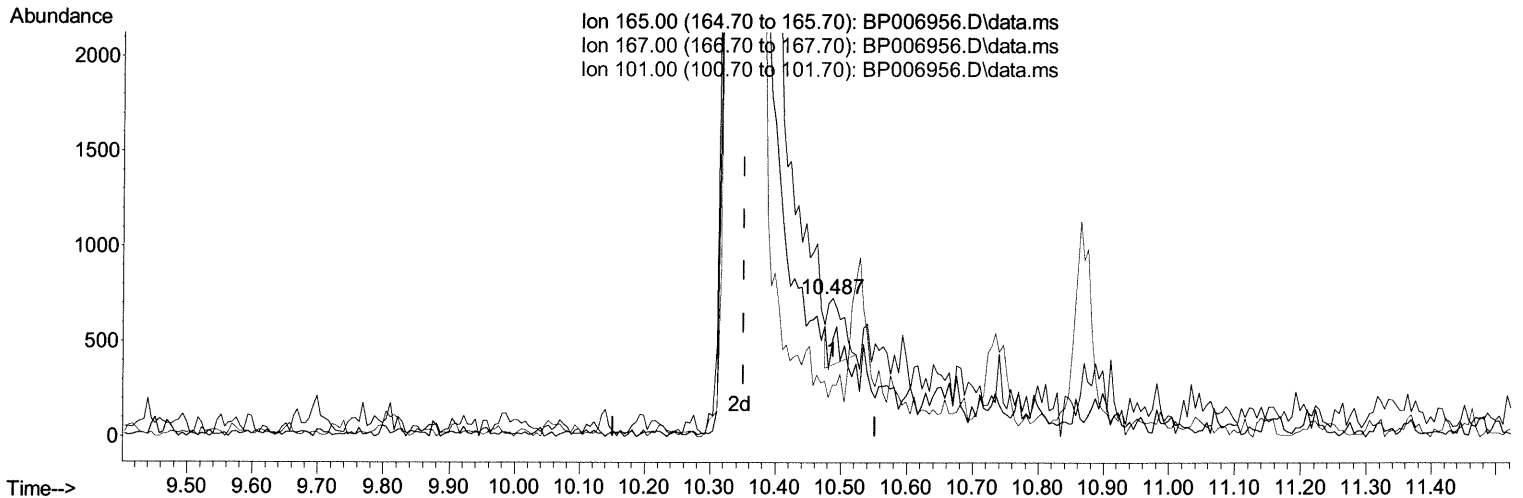
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(28) 2,4-Dichlorophenol-d3 (S)

10.487min (+ 0.135) 0.03 ng/ul

response 552

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	62.40	70.37
101.00	40.60	37.59
0.00	0.00	0.00

Quantitation Report (Qedit)

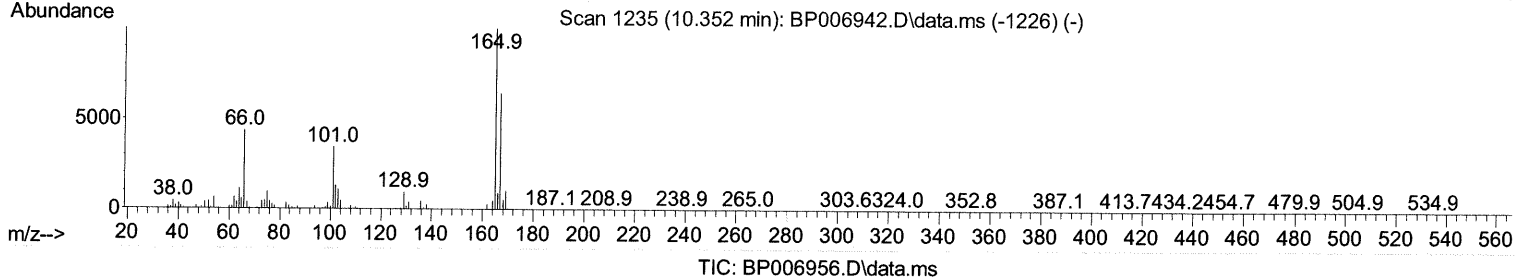
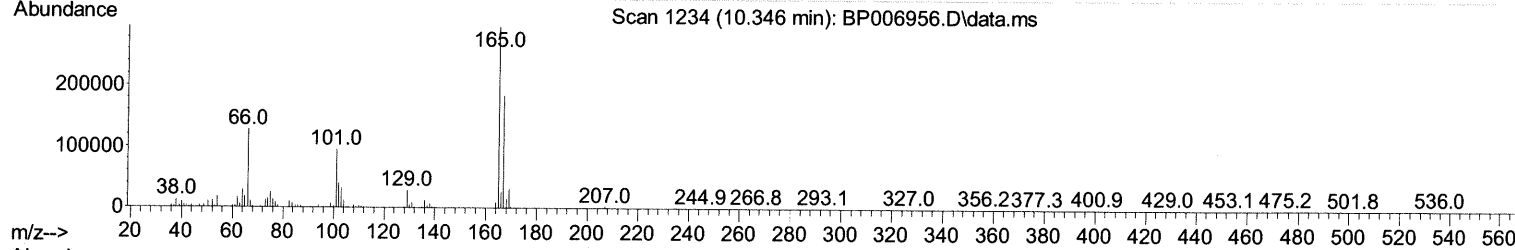
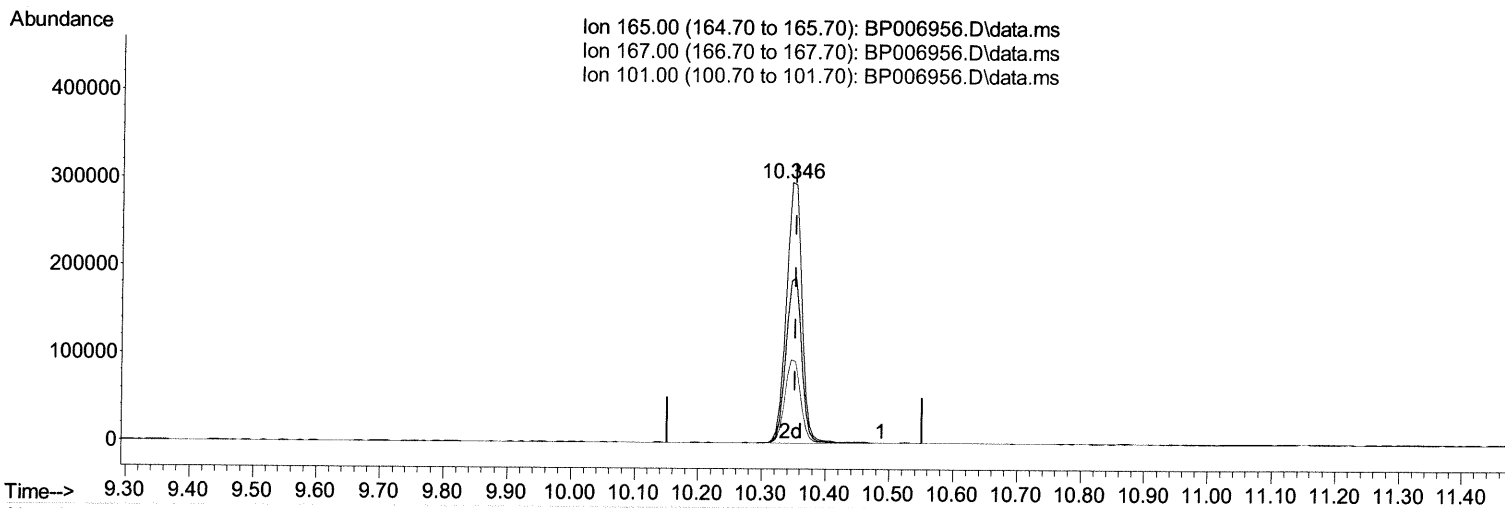
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(28) 2,4-Dichlorophenol-d3 (S)

10.346min (-0.006) 31.67 ng/ul m 09/14/21 34

response 499794

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	62.40	62.20
101.00	40.60	32.06#
0.00	0.00	0.00

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 Sample : M3651-12
 Misc :
 ALS Vial : 56 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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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	275465	20.000	ng/ul	0.00
20) Naphthalene-d8	10.734	136	1122317	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.557	164	700290	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.304	188	1471607	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.386	240	1484149	20.000	ng/ul	#-0.01
88) Perylene-d12	23.816	264	1537833	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	17210	2.486	ng/ul	0.00
4) Pyridine-d5	3.793	84	170030	9.426	ng/ul	0.01
7) Phenol-d5	7.087	99	169924	8.393	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	416170	35.791	ng/ul	0.00
11) 2-Chlorophenol-d4	7.463	132	465803	28.499	ng/ul	0.00
15) 4-Methylphenol-d8	8.634	113	299007	18.807	ng/ul	0.00
21) Nitrobenzene-d5	9.093	128	257701	37.908	ng/ul	0.00
24) 2-Nitrophenol-d4	9.816	143	254696	38.056	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.346	165	499794m	31.671	ng/ul	0.00
31) 4-Chloroaniline-d4	10.869	131	640793	29.362	ng/ul	0.00
46) Dimethylphthalate-d6	13.969	166	1687373	37.203	ng/ul	0.00
49) Acenaphthylene-d8	14.251	160	2037359	35.467	ng/ul	0.00
54) 4-Nitrophenol-d4	14.740	143	44103	5.597	ng/ul	0.00
60) Fluorene-d10	15.545	176	1543158	38.550	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.657	200	197782	30.381	ng/ul	-0.01
73) Anthracene-d10	17.404	188	2649444	43.487	ng/ul	-0.01
81) Pyrene-d10	19.633	212	3126470	44.981	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.657	264	3087326	41.614	ng/ul	-0.02

Target Compounds Qvalue

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