

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070823\
 Data File : BP016140.D
 Acq On : 09 Jul 2023 19:37
 Operator : MA/JU
 Sample : 03356-06RE
 Misc :
 ALS Vial : 65 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Jul 09 22:44:39 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jul 08 10:11:43 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.016	152	341966	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.846	136	1450090	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.669	164	839724	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.439	188	1622277	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.551	240	1214130	20.000	ng/ul	# 0.01
88) Perylene-d12	24.080	264	1298806	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	32938	3.465	ng/uL	0.00
4) Pyridine-d5	3.799	84	388397	16.211	ng/ul	0.00
7) Phenol-d5	7.222	99	510607	17.451	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.352	67	288009	15.910	ng/ul	0.00
11) 2-Chlorophenol-d4	7.558	132	403087	18.250	ng/ul	0.00
15) 4-Methylphenol-d8	8.775	113	416372	18.014	ng/ul	0.00
21) Nitrobenzene-d5	9.228	128	195357	17.628	ng/ul	0.01
24) 2-Nitrophenol-d4	9.969	143	217264	16.798	ng/ul	0.02
28) 2,4-Dichlorophenol-d3	10.563	165	287175	12.459	ng/ul	0.06
31) 4-Chloroaniline-d4	11.063	131	428815	14.483	ng/ul	0.04
46) Dimethylphthalate-d6	14.098	166	1222581	18.837	ng/ul	0.01
49) Acenaphthylene-d8	14.369	160	1386414	19.002	ng/ul	0.00
54) 4-Nitrophenol-d4	15.075	143	29222	3.412	ng/ul	0.11
60) Fluorene-d10	15.681	176	1026524	19.208	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.910	200	66624	6.678	ng/ul	0.06
73) Anthracene-d10	17.551	188	1476477	19.925	ng/ul	0.00
81) Pyrene-d10	19.775	212	1582441	19.960	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.904	264	1225801	18.709	ng/ul	0.00

Target Compounds	Qvalue

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

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