

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110123\
 Data File : BP017854.D
 Acq On : 01 Nov 2023 16:40
 Operator : MA/JU
 Sample : 04952-12
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EOAS6

Quant Time: Nov 02 01:13:56 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 19 01:32:15 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.881	152	132613	20.000	ng/ul	0.00
20) Naphthalene-d8	10.710	136	493288	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.575	164	284595	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.475	188	1163596	20.000	ng/ul	0.10
79) Chrysene-d12	21.516	240	466834	20.000	ng/ul	0.00
88) Perylene-d12	24.074	264	503821	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	11837	3.321	ng/uL	0.01
4) Pyridine-d5	3.711	84	105091	11.260	ng/ul	0.02
7) Phenol-d5	7.046	99	64938	5.885	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	238138	34.942	ng/ul	0.00
11) 2-Chlorophenol-d4	7.404	132	220949	24.548	ng/ul	0.00
15) 4-Methylphenol-d8	8.604	113	130871	15.102	ng/ul	0.00
21) Nitrobenzene-d5	9.087	128	149740	37.462	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.798	143	154650	35.017	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.334	165	244504	28.223	ng/ul	0.00
31) 4-Chloroaniline-d4	10.887	131	324030	28.580	ng/ul	0.00
46) Dimethylphthalate-d6	13.992	166	879498	36.813	ng/ul	-0.01
49) Acenaphthylene-d8	14.275	160	966902	36.291	ng/ul	0.00
54) 4-Nitrophenol-d4	14.828	143	98	0.029	ng/ul	0.00
60) Fluorene-d10	15.581	176	731573	35.424	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.910	200	138	0.018	ng/ul	0.18
73) Anthracene-d10	17.475	188	1166214	19.613	ng/ul	0.00
81) Pyrene-d10	19.733	212	1315973	44.963	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.910	264	1104142	39.347	ng/ul	0.01

Target Compounds Qvalue

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

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