

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090723\
 Data File : BP016994.D
 Acq On : 07 Sep 2023 18:45
 Operator : MA/JU
 Sample : 04123-11
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH9H3

Quant Time: Sep 08 01:28:29 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 06 01:52:05 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.975	152	340742	20.000	ng/u1	0.00
20) Naphthalene-d8	10.805	136	1512069	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.634	164	928236	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.404	188	2054521	20.000	ng/u1	0.00
79) Chrysene-d12	21.504	240	1818873	20.000	ng/u1	0.00
88) Perylene-d12	24.045	264	1855140	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.323	96	64549	7.166	ng/uL	0.00
4) Pyridine-d5	3.764	84	326648	12.690	ng/u1	0.00
7) Phenol-d5	7.122	99	306960	9.861	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.317	67	805802	41.383	ng/u1	0.00
11) 2-Chlorophenol-d4	7.499	132	742215	32.943	ng/u1	0.00
15) 4-Methylphenol-d8	8.687	113	543363	22.911	ng/u1	0.00
21) Nitrobenzene-d5	9.175	128	473231	40.341	ng/u1	0.00
24) 2-Nitrophenol-d4	9.893	143	499929	39.314	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.416	165	838022	35.696	ng/u1	0.00
31) 4-Chloroaniline-d4	10.963	131	1050166	31.005	ng/u1	0.00
46) Dimethylphthalate-d6	14.051	166	2938307	42.525	ng/u1	0.00
49) Acenaphthylene-d8	14.340	160	3341358	41.903	ng/u1	0.00
54) 4-Nitrophenol-d4	14.851	143	118032	8.745	ng/u1	0.00
60) Fluorene-d10	15.634	176	2557853	43.237	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.757	200	438764	35.487	ng/u1	0.00
73) Anthracene-d10	17.504	188	4103580	43.492	ng/u1	0.00
81) Pyrene-d10	19.745	212	4811477	43.430	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.874	264	4354066	46.263	ng/u1	0.00

Target Compounds Qvalue

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

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