

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
 Data File : BP008538.D
 Acq On : 23 Dec 2021 22:46
 Operator : CG/JU
 Sample : M5090-15
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BG3C7

Quant Time: Dec 24 04:14:42 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP122321.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 23 14:48:11 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.904	152	503508	20.000	ng/u1	0.00
20) Naphthalene-d8	10.704	136	2181957	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.534	164	1523920	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.280	188	3335350	20.000	ng/u1	0.00
79) Chrysene-d12	21.363	240	3460186	20.000	ng/u1	0.00
88) Perylene-d12	23.792	264	3520957	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	52244	4.677	ng/uL	0.00
4) Pyridine-d5	3.758	84	306002	9.789	ng/u1	0.00
7) Phenol-d5	7.057	99	260718	6.491	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	685357	30.041	ng/u1	0.00
11) 2-Chlorophenol-d4	7.434	132	767149	24.008	ng/u1	0.00
15) 4-Methylphenol-d8	8.604	113	526428	16.337	ng/u1	0.00
21) Nitrobenzene-d5	9.057	128	487453	31.571	ng/u1	0.00
24) 2-Nitrophenol-d4	9.781	143	458013	32.353	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.322	165	1015164	29.213	ng/u1	0.00
31) 4-Chloroaniline-d4	10.834	131	1226229	25.563	ng/u1	0.00
46) Dimethylphthalate-d6	13.945	166	3989666	35.433	ng/u1	0.00
49) Acenaphthylene-d8	14.228	160	4673439	33.194	ng/u1	0.00
54) 4-Nitrophenol-d4	14.710	143	111559	6.250	ng/u1	0.00
60) Fluorene-d10	15.528	176	3436410	35.152	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.633	200	415802	28.320	ng/u1	0.00
73) Anthracene-d10	17.380	188	5775445	36.541	ng/u1	0.00
81) Pyrene-d10	19.610	212	6834537	32.963	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.639	264	6769019	37.353	ng/u1	0.00

Target Compounds Qvalue

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

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