

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122721\
 Data File : BP008548.D
 Acq On : 27 Dec 2021 10:37
 Operator : CG/JU
 Sample : PB141660BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK660

Quant Time: Dec 27 11:14:08 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	7.905	152	527447	20.000	ng/u1	0.00
20) Naphthalene-d8	10.704	136	2092092	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.534	164	1401232	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.281	188	2840476	20.000	ng/u1	0.00
79) Chrysene-d12	21.363	240	2163197	20.000	ng/u1	0.00
88) Perylene-d12	23.792	264	1994949	20.000	ng/u1	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	71676	6.125	ng/uL	0.00
4) Pyridine-d5	3.758	84	921665	28.145	ng/u1	0.00
7) Phenol-d5	7.058	99	1161111	27.595	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	661973	27.699	ng/u1	0.00
11) 2-Chlorophenol-d4	7.434	132	964532	28.816	ng/u1	0.00
15) 4-Methylphenol-d8	8.604	113	912603	27.036	ng/u1	0.00
21) Nitrobenzene-d5	9.057	128	452365	30.557	ng/u1	0.00
24) 2-Nitrophenol-d4	9.781	143	428928	31.600	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.322	165	933312	28.011	ng/u1	0.00
31) 4-Chloroaniline-d4	10.834	131	1346714	29.280	ng/u1	0.00
46) Dimethylphthalate-d6	13.945	166	3130673	30.239	ng/u1	0.00
49) Acenaphthylene-d8	14.228	160	3800038	29.354	ng/u1	0.00
54) 4-Nitrophenol-d4	14.710	143	395197	24.081	ng/u1	0.00
60) Fluorene-d10	15.528	176	2684678	29.867	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.634	200	264753	21.174	ng/u1	0.00
73) Anthracene-d10	17.381	188	4205190	31.241	ng/u1	0.00
81) Pyrene-d10	19.610	212	4447115	34.308	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.633	264	3208792	31.252	ng/u1	0.00

Target Compounds	Qvalue

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

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