

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122721\
 Data File : BP008572.D
 Acq On : 28 Dec 2021 09:39
 Operator : CG/JU
 Sample : PB141660BL
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK660

Quant Time: Dec 28 12:14:58 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	550906	20.000	ng/u1	0.00
20) Naphthalene-d8	10.710	136	2285847	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.540	164	1668162	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.287	188	3706538	20.000	ng/u1	0.00
79) Chrysene-d12	21.369	240	3831947	20.000	ng/u1	0.00
88) Perylene-d12	23.810	264	3655395	20.000	ng/u1	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	79092	6.471	ng/uL	0.00
4) Pyridine-d5	3.758	84	728757	21.306	ng/u1	0.00
7) Phenol-d5	7.064	99	1316482	29.955	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.234	67	749132	30.011	ng/u1	0.00
11) 2-Chlorophenol-d4	7.434	132	1095402	31.332	ng/u1	0.00
15) 4-Methylphenol-d8	8.611	113	1072858	30.430	ng/u1	0.00
21) Nitrobenzene-d5	9.063	128	548929	33.937	ng/u1	0.00
24) 2-Nitrophenol-d4	9.787	143	537574	36.247	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.322	165	1149249	31.568	ng/u1	0.00
31) 4-Chloroaniline-d4	10.840	131	738328	14.692	ng/u1	0.00
46) Dimethylphthalate-d6	13.951	166	4280968	34.733	ng/u1	0.00
49) Acenaphthylene-d8	14.234	160	4952212	32.133	ng/u1	0.00
54) 4-Nitrophenol-d4	14.722	143	661536	33.860	ng/u1	0.00
60) Fluorene-d10	15.534	176	3633743	33.957	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.640	200	466203	28.573	ng/u1	0.00
73) Anthracene-d10	17.387	188	6069226	34.554	ng/u1	0.00
81) Pyrene-d10	19.616	212	7123671	31.024	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.651	264	6576502	34.956	ng/u1	0.01

Target Compounds Qvalue

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

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