

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080322\
 Data File : BP011241.D
 Acq On : 03 Aug 2022 12:55
 Operator : CG/JU
 Sample : PB146573BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK573

Quant Time: Aug 03 22:59:36 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080222.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 02 23:52:37 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.599	152	183527	20.000	ng/u1	0.00
20) Naphthalene-d8	10.393	136	830765	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.275	164	442080	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.045	188	859022	20.000	ng/u1	0.00
79) Chrysene-d12	21.175	240	847037	20.000	ng/u1	0.00
88) Perylene-d12	23.498	264	920681	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.093	96	9904	4.363	ng/uL	0.00
4) Pyridine-d5	3.505	84	152690	24.939	ng/u1	0.00
7) Phenol-d5	6.781	99	416355	27.223	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.946	67	321045	29.090	ng/u1	0.00
11) 2-Chlorophenol-d4	7.134	132	333894	26.783	ng/u1	0.00
15) 4-Methylphenol-d8	8.322	113	324779	25.807	ng/u1	0.00
21) Nitrobenzene-d5	8.763	128	172083	28.798	ng/u1	0.00
24) 2-Nitrophenol-d4	9.487	143	135214	24.140	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.022	165	274818	25.729	ng/u1	0.00
31) 4-Chloroaniline-d4	10.546	131	451212	27.306	ng/u1	0.00
46) Dimethylphthalate-d6	13.699	166	892538	29.478	ng/u1	0.00
49) Acenaphthylene-d8	13.963	160	1153039	30.267	ng/u1	0.00
54) 4-Nitrophenol-d4	14.504	143	83981	16.311	ng/u1	0.00
60) Fluorene-d10	15.281	176	802425	31.146	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.410	200	63759	15.985	ng/u1	0.00
73) Anthracene-d10	17.145	188	1204786	31.912	ng/u1	0.00
81) Pyrene-d10	19.410	212	1318819	30.794	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.345	264	1378174	30.039	ng/u1	0.00

Target Compounds Qvalue

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

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