

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080322\
 Data File : BP011237.D
 Acq On : 03 Aug 2022 09:47
 Operator : CG/JU
 Sample : PB146573BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK573

Quant Time: Aug 03 22:58:31 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	189372	20.000	ng/u1	0.00
20) Naphthalene-d8	10.393	136	862827	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.275	164	468267	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.045	188	895545	20.000	ng/u1	0.00
79) Chrysene-d12	21.174	240	841411	20.000	ng/u1	0.00
88) Perylene-d12	23.492	264	919928	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.093	96	10455	4.464	ng/uL	0.00
4) Pyridine-d5	3.505	84	168574	26.684	ng/u1	0.00
7) Phenol-d5	6.781	99	436144	27.636	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.946	67	327916	28.796	ng/u1	0.00
11) 2-Chlorophenol-d4	7.134	132	351624	27.335	ng/u1	0.00
15) 4-Methylphenol-d8	8.322	113	352621	27.155	ng/u1	0.00
21) Nitrobenzene-d5	8.763	128	178214	28.716	ng/u1	0.00
24) 2-Nitrophenol-d4	9.487	143	150857	25.932	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.022	165	293296	26.439	ng/u1	0.00
31) 4-Chloroaniline-d4	10.546	131	475962	27.734	ng/u1	0.00
46) Dimethylphthalate-d6	13.698	166	964003	30.058	ng/u1	0.00
49) Acenaphthylene-d8	13.963	160	1222919	30.306	ng/u1	0.00
54) 4-Nitrophenol-d4	14.498	143	93081	17.067	ng/u1	0.00
60) Fluorene-d10	15.281	176	827137	30.310	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.410	200	68455	16.462	ng/u1	0.00
73) Anthracene-d10	17.145	188	1221639	31.039	ng/u1	0.00
81) Pyrene-d10	19.404	212	1318411	30.991	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.345	264	1358794	29.641	ng/u1	0.00

Target Compounds	Qvalue

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

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