

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080122\
 Data File : BP011194.D
 Acq On : 01 Aug 2022 11:54
 Operator : CG/JU
 Sample : PB146474BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK474

Quant Time: Aug 01 22:59:11 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP072822.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jul 28 16:05:22 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.652	152	308561	20.000	ng/u1	0.00
20) Naphthalene-d8	10.446	136	1385475	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.316	164	691454	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.086	188	1148169	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.216	240	1030356	20.000	ng/u1	# 0.00
88) Perylene-d12	23.563	264	1104690	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.152	96	24808	3.682	ng/uL	0.00
4) Pyridine-d5	3.558	84	414175	21.728	ng/u1	0.00
7) Phenol-d5	6.828	99	771290	28.638	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.993	67	542094	28.465	ng/u1	0.00
11) 2-Chlorophenol-d4	7.181	132	601303	28.331	ng/u1	0.00
15) 4-Methylphenol-d8	8.369	113	610393	27.705	ng/u1	0.00
21) Nitrobenzene-d5	8.816	128	310061	28.796	ng/u1	0.00
24) 2-Nitrophenol-d4	9.534	143	295603	28.661	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.069	165	487377	26.814	ng/u1	0.00
31) 4-Chloroaniline-d4	10.593	131	800769	27.176	ng/u1	0.00
46) Dimethylphthalate-d6	13.740	166	1400577	29.730	ng/u1	0.00
49) Acenaphthylene-d8	14.010	160	1766020	29.831	ng/u1	0.00
54) 4-Nitrophenol-d4	14.540	143	148160	17.484	ng/u1	0.00
60) Fluorene-d10	15.322	176	1116981	28.217	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.451	200	105788	18.278	ng/u1	0.00
73) Anthracene-d10	17.186	188	1502687	29.702	ng/u1	0.00
81) Pyrene-d10	19.445	212	1576756	28.479	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.410	264	1637856	29.850	ng/u1	0.00

Target Compounds Qvalue

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

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