

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080122\
 Data File : BP011211.D
 Acq On : 01 Aug 2022 22:12
 Operator : CG/JU
 Sample : PB146356BL
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK356

Quant Time: Aug 02 00:29:42 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	7.646	152	288278	20.000	ng/ul	0.00
20) Naphthalene-d8	10.440	136	1215469	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.316	164	612778	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.087	188	1156024	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.210	240	1134363	20.000	ng/ul	# 0.00
88) Perylene-d12	23.551	264	1191629	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.146	96	23909	3.798	ng/uL	0.00
4) Pyridine-d5	3.558	84	351840	19.757	ng/ul	0.00
7) Phenol-d5	6.828	99	750568	29.829	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.993	67	526781	29.607	ng/ul	0.00
11) 2-Chlorophenol-d4	7.181	132	611340	30.830	ng/ul	0.00
15) 4-Methylphenol-d8	8.364	113	574713	27.921	ng/ul	0.00
21) Nitrobenzene-d5	8.811	128	289628	30.661	ng/ul	0.00
24) 2-Nitrophenol-d4	9.528	143	288757	31.913	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.063	165	468731	29.395	ng/ul	0.00
31) 4-Chloroaniline-d4	10.587	131	729717	28.228	ng/ul	0.00
46) Dimethylphthalate-d6	13.734	166	1340791	32.115	ng/ul	0.00
49) Acenaphthylene-d8	14.004	160	1697357	32.353	ng/ul	0.00
54) 4-Nitrophenol-d4	14.540	143	178567	23.778	ng/ul	0.00
60) Fluorene-d10	15.316	176	1103018	31.441	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.451	200	116045	19.913	ng/ul	0.00
73) Anthracene-d10	17.187	188	1643974	32.274	ng/ul	0.00
81) Pyrene-d10	19.445	212	1857411	30.472	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.404	264	1908256	32.241	ng/ul	0.00

Target Compounds Qvalue

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

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