

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080422\
 Data File : BP011258.D
 Acq On : 04 Aug 2022 11:25
 Operator : CG/JU
 Sample : PB146590BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK590

Quant Time: Aug 04 22:53:10 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	167055	20.000	ng/u1	0.00
20) Naphthalene-d8	10.393	136	736460	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.275	164	427463	20.000	ng/u1	# 0.00
64) Phenanthrene-d10	17.045	188	912682	20.000	ng/u1	0.00
79) Chrysene-d12	21.174	240	1022285	20.000	ng/u1	0.00
88) Perylene-d12	23.492	264	1112775	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.093	96	7205	3.487	ng/uL	0.00
4) Pyridine-d5	3.505	84	121337	21.772	ng/u1	0.00
7) Phenol-d5	6.781	99	357065	25.648	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.946	67	299279	29.792	ng/u1	0.00
11) 2-Chlorophenol-d4	7.134	132	287515	25.337	ng/u1	0.00
15) 4-Methylphenol-d8	8.322	113	276721	24.157	ng/u1	0.00
21) Nitrobenzene-d5	8.763	128	166135	31.363	ng/u1	0.00
24) 2-Nitrophenol-d4	9.487	143	120902	24.349	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.022	165	235260	24.846	ng/u1	0.00
31) 4-Chloroaniline-d4	10.546	131	372328	25.417	ng/u1	0.00
46) Dimethylphthalate-d6	13.698	166	842878	28.790	ng/u1	0.00
49) Acenaphthylene-d8	13.963	160	1140109	30.951	ng/u1	0.00
54) 4-Nitrophenol-d4	14.504	143	94343	18.950	ng/u1	0.00
60) Fluorene-d10	15.275	176	850393	34.137	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.410	200	85754	20.235	ng/u1	0.00
73) Anthracene-d10	17.145	188	1494583	37.261	ng/u1	0.00
81) Pyrene-d10	19.410	212	1690001	32.697	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.345	264	1829771	32.998	ng/u1	0.00

Target Compounds	Qvalue

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

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