

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
 Data File : BP012208.D
 Acq On : 20 Oct 2022 09:59
 Operator : CG/JU
 Sample : PB148325BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK325

Quant Time: Oct 20 23:21:15 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 20 00:51:41 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.822	152	364526	20.000	ng/u1	0.00
20) Naphthalene-d8	10.628	136	1573221	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.475	164	1041396	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.234	188	2258721	20.000	ng/u1	0.00
79) Chrysene-d12	21.327	240	1978185	20.000	ng/u1	0.00
88) Perylene-d12	23.733	264	1674479	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	52354	5.812	ng/uL	0.00
4) Pyridine-d5	3.681	84	628334	24.209	ng/u1	0.00
7) Phenol-d5	6.993	99	893795	28.401	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.158	67	558929	29.286	ng/u1	0.00
11) 2-Chlorophenol-d4	7.358	132	713919	29.803	ng/u1	0.00
15) 4-Methylphenol-d8	8.540	113	717431	29.332	ng/u1	0.00
21) Nitrobenzene-d5	8.993	128	355779	30.552	ng/u1	0.00
24) 2-Nitrophenol-d4	9.716	143	391047	31.124	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.252	165	683672	29.010	ng/u1	0.00
31) 4-Chloroaniline-d4	10.775	131	966456	28.863	ng/u1	0.00
46) Dimethylphthalate-d6	13.893	166	2366285	32.665	ng/u1	0.00
49) Acenaphthylene-d8	14.169	160	2555851	30.910	ng/u1	0.00
54) 4-Nitrophenol-d4	14.687	143	372529	27.449	ng/u1	0.00
60) Fluorene-d10	15.475	176	2015405	32.339	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	326218	25.387	ng/u1	0.00
73) Anthracene-d10	17.334	188	3272837	33.057	ng/u1	0.00
81) Pyrene-d10	19.575	212	3650356	32.411	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.580	264	2872556	34.208	ng/u1	0.00

Target Compounds Qvalue

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

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