

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
 Data File : BP012209.D
 Acq On : 20 Oct 2022 10:33
 Operator : CG/JU
 Sample : PB148327BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK327

Quant Time: Oct 20 23:21:25 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	7.828	152	387413	20.000	ng/u1	0.00
20) Naphthalene-d8	10.628	136	1703520	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.475	164	1099711	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.234	188	2367977	20.000	ng/u1	0.00
79) Chrysene-d12	21.328	240	2222109	20.000	ng/u1	0.00
88) Perylene-d12	23.733	264	2003536	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	60095	6.278	ng/uL	0.00
4) Pyridine-d5	3.681	84	748818	27.146	ng/u1	0.00
7) Phenol-d5	6.993	99	1037433	31.018	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.158	67	662308	32.652	ng/u1	0.00
11) 2-Chlorophenol-d4	7.358	132	820987	32.248	ng/u1	0.00
15) 4-Methylphenol-d8	8.540	113	817694	31.457	ng/u1	0.00
21) Nitrobenzene-d5	8.993	128	412936	32.748	ng/u1	0.00
24) 2-Nitrophenol-d4	9.716	143	450721	33.130	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.252	165	770287	30.185	ng/u1	0.00
31) 4-Chloroaniline-d4	10.775	131	1114190	30.730	ng/u1	0.00
46) Dimethylphthalate-d6	13.893	166	2690535	35.172	ng/u1	0.00
49) Acenaphthylene-d8	14.169	160	2924341	33.491	ng/u1	0.00
54) 4-Nitrophenol-d4	14.687	143	438800	30.618	ng/u1	0.00
60) Fluorene-d10	15.475	176	2281238	34.663	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	401188	29.781	ng/u1	0.00
73) Anthracene-d10	17.334	188	3698518	35.633	ng/u1	0.00
81) Pyrene-d10	19.575	212	4193341	33.145	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.580	264	3686201	36.688	ng/u1	0.00

Target Compounds Qvalue

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

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