

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042922\
 Data File : BP010160.D
 Acq On : 29 Apr 2022 20:38
 Operator : CG/JU
 Sample : PB144447BL
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK447

Quant Time: Apr 30 02:38:26 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 21 14:49:38 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	111552	20.000	ng/u1	-0.02
20) Naphthalene-d8	10.728	136	425680	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.557	164	249443	20.000	ng/u1	-0.02
64) Phenanthrene-d10	17.310	188	503940	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.398	240	486141	20.000	ng/u1	#-0.01
88) Perylene-d12	23.845	264	496909	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.357	96	20509	8.017	ng/uL	-0.01
4) Pyridine-d5	3.775	84	253080	35.167	ng/u1	-0.01
7) Phenol-d5	7.087	99	330201	32.543	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.251	67	224861	36.932	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.457	132	279777	37.231	ng/u1	-0.01
15) 4-Methylphenol-d8	8.634	113	267134	31.776	ng/u1	-0.02
21) Nitrobenzene-d5	9.081	128	133838	39.531	ng/u1	-0.02
24) 2-Nitrophenol-d4	9.810	143	146612	39.651	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.351	165	243448	33.821	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.863	131	341687	33.106	ng/u1	-0.02
46) Dimethylphthalate-d6	13.969	166	774517	39.916	ng/u1	-0.01
49) Acenaphthylene-d8	14.251	160	890933	37.938	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.757	143	104409	29.228	ng/u1	0.00
60) Fluorene-d10	15.551	176	654067	39.186	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.669	200	99423	31.201	ng/u1	-0.01
73) Anthracene-d10	17.410	188	984055	40.363	ng/u1	-0.01
81) Pyrene-d10	19.645	212	1153776	41.732	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.686	264	1101365	42.278	ng/u1	-0.01

Target Compounds Qvalue

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

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