

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042023\
 Data File : BP014809.D
 Acq On : 22 Apr 2023 01:52
 Operator : CG/JU
 Sample : PB152279BL
 Misc :
 ALS Vial : 64 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK279

Quant Time: Apr 22 04:07:18 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041823.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Apr 22 04:03:47 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.169	152	309109	20.000	ng/u1	0.02
20) Naphthalene-d8	10.999	136	1312162	20.000	ng/u1	0.02
38) Acenaphthene-d10	14.798	164	842230	20.000	ng/u1	0.01
64) Phenanthrene-d10	17.545	188	1794470	20.000	ng/u1	0.01
79) Chrysene-d12	21.616	240	1277616	20.000	ng/u1	0.01
88) Perylene-d12	24.204	264	1128956	20.000	ng/u1	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.499	96	51418	6.611	ng/uL	0.01
4) Pyridine-d5	3.928	84	534837	25.258	ng/u1	0.01
7) Phenol-d5	7.311	99	868667	32.503	ng/u1	0.02
9) Bis-(2-Chloroethyl)eth...	7.487	67	548478	33.265	ng/u1	0.02
11) 2-Chlorophenol-d4	7.693	132	700862	35.311	ng/u1	0.02
15) 4-Methylphenol-d8	8.869	113	720357	34.574	ng/u1	0.02
21) Nitrobenzene-d5	9.346	128	331694	40.341	ng/u1	0.02
24) 2-Nitrophenol-d4	10.069	143	336023	45.248	ng/u1	0.02
28) 2,4-Dichlorophenol-d3	10.610	165	675242	33.487	ng/u1	0.02
31) 4-Chloroaniline-d4	11.128	131	500614	16.266	ng/u1	0.02
46) Dimethylphthalate-d6	14.198	166	2361271	35.407	ng/u1	0.02
49) Acenaphthylene-d8	14.493	160	2654444	34.346	ng/u1	0.01
54) 4-Nitrophenol-d4	14.969	143	335291	31.609	ng/u1	0.02
60) Fluorene-d10	15.787	176	2003132	35.006	ng/u1	0.01
65) 4,6-Dinitro-2-methylph...	15.893	200	245390	33.791	ng/u1	0.01
73) Anthracene-d10	17.645	188	3094425	35.696	ng/u1	0.01
81) Pyrene-d10	19.857	212	3289436	44.945	ng/u1	0.01
92) Benzo(a)pyrene-d12	24.033	264	2205981	37.701	ng/u1	0.02

Target Compounds Qvalue

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

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