

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
 Data File : BM044981.D
 Acq On : 06 Apr 2024 22:03
 Operator : MA/JU
 Sample : PB160053BL
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK053

Quant Time: Apr 08 05:04:27 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Apr 08 05:01:54 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.634	152	79197	20.000	ng/u1	0.00
20) Naphthalene-d8	10.410	136	350046	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.280	164	227359	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.039	188	490963	20.000	ng/u1	0.00
79) Chrysene-d12	21.250	240	462895	20.000	ng/u1	0.00
88) Perylene-d12	23.474	264	557060	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.210	96	11920	5.620	ng/uL	0.00
4) Pyridine-d5	3.610	84	177844	30.836	ng/u1	0.00
7) Phenol-d5	6.816	99	214519	31.951	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.987	67	130789	32.697	ng/u1	0.00
11) 2-Chlorophenol-d4	7.169	132	179429	34.112	ng/u1	0.00
15) 4-Methylphenol-d8	8.345	113	174098	33.128	ng/u1	0.00
21) Nitrobenzene-d5	8.798	128	88260	32.825	ng/u1	0.00
24) 2-Nitrophenol-d4	9.504	143	99647	35.112	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.039	165	168119	32.310	ng/u1	0.00
31) 4-Chloroaniline-d4	10.563	131	289263	34.917	ng/u1	0.00
46) Dimethylphthalate-d6	13.710	166	513902	32.463	ng/u1	0.00
49) Acenaphthylene-d8	13.974	160	615001	32.885	ng/u1	0.00
54) 4-Nitrophenol-d4	14.521	143	88955	26.925	ng/u1	0.00
60) Fluorene-d10	15.286	176	461645	32.710	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.421	200	75209	25.836	ng/u1	0.00
73) Anthracene-d10	17.139	188	745828	33.769	ng/u1	0.00
81) Pyrene-d10	19.445	212	865786	37.985	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.333	264	948252	34.894	ng/u1	0.00

Target Compounds Qvalue

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

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