

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
 Data File : BM044965.D
 Acq On : 06 Apr 2024 11:52
 Operator : MA/JU
 Sample : PB160048BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Apr 08 01:49:40 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Apr 08 01:45:36 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.634	152	72399	20.000	ng/ul	0.00
20) Naphthalene-d8	10.410	136	308808	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.280	164	193915	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.039	188	400555	20.000	ng/ul	0.00
79) Chrysene-d12	21.250	240	362091	20.000	ng/ul	0.00
88) Perylene-d12	23.474	264	447199	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.210	96	10428	5.379	ng/uL	0.00
4) Pyridine-d5	3.610	84	115035	21.819	ng/ul	0.00
7) Phenol-d5	6.816	99	138220	22.520	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.987	67	89134	24.375	ng/ul	0.00
11) 2-Chlorophenol-d4	7.169	132	119405	24.832	ng/ul	0.00
15) 4-Methylphenol-d8	8.340	113	111132	23.132	ng/ul	0.00
21) Nitrobenzene-d5	8.798	128	56291	23.731	ng/ul	0.00
24) 2-Nitrophenol-d4	9.510	143	59402	23.726	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.039	165	102238	22.272	ng/ul	0.00
31) 4-Chloroaniline-d4	10.569	131	151188	20.687	ng/ul	0.00
46) Dimethylphthalate-d6	13.710	166	322562	23.890	ng/ul	0.00
49) Acenaphthylene-d8	13.974	160	382143	23.958	ng/ul	0.00
54) 4-Nitrophenol-d4	14.521	143	42022	14.913	ng/ul	0.00
60) Fluorene-d10	15.286	176	289778	24.074	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.427	200	38259	16.109	ng/ul	0.00
73) Anthracene-d10	17.145	188	448961	24.916	ng/ul	0.00
81) Pyrene-d10	19.445	212	502055	28.159	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.333	264	544000	24.936	ng/ul	0.00

Target Compounds Qvalue

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

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