

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080624\
 Data File : BM047067.D
 Acq On : 07 Aug 2024 21:28
 Operator : RC/JU
 Sample : PB162513TB
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLEB466

Quant Time: Aug 08 00:12:29 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Aug 07 04:48:25 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.392	152	295716	20.000	ng/u1	0.00
20) Naphthalene-d8	10.151	136	1115857	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.045	164	698952	20.000	ng/u1	0.00
64) Phenanthrene-d10	16.804	188	1408648	20.000	ng/u1	-0.01
79) Chrysene-d12	21.039	240	1207946	20.000	ng/u1	-0.01
88) Perylene-d12	23.744	264	1345563	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.016	96	36650	4.910	ng/uL	0.00
4) Pyridine-d5	3.404	84	535636	26.308	ng/u1	0.00
7) Phenol-d5	6.598	99	649570	28.916	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.751	67	413537	27.978	ng/u1	0.00
11) 2-Chlorophenol-d4	6.934	132	515648	27.098	ng/u1	0.00
15) 4-Methylphenol-d8	8.110	113	499695	28.080	ng/u1	-0.01
21) Nitrobenzene-d5	8.539	128	258731	28.895	ng/u1	0.00
24) 2-Nitrophenol-d4	9.251	143	285052	29.151	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	9.786	165	511335	27.236	ng/u1	0.00
31) 4-Chloroaniline-d4	10.298	131	688529	27.905	ng/u1	0.00
46) Dimethylphthalate-d6	13.474	166	1476709	27.632	ng/u1	0.00
49) Acenaphthylene-d8	13.733	160	1697953	28.463	ng/u1	0.00
54) 4-Nitrophenol-d4	14.286	143	253111	24.495	ng/u1	-0.01
60) Fluorene-d10	15.051	176	1239982	27.123	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.192	200	177310	19.923	ng/u1	0.00
73) Anthracene-d10	16.904	188	1912467	28.639	ng/u1	0.00
81) Pyrene-d10	19.215	212	2076279	30.026	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.562	264	2036061	28.736	ng/u1	-0.01

Target Compounds Qvalue

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

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