

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080624\
 Data File : BM047042.D
 Acq On : 07 Aug 2024 04:42
 Operator : RC/JU
 Sample : PB162282BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK282

Quant Time: Aug 07 08:40:35 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	7.398	152	233089	20.000	ng/u1	0.00
20) Naphthalene-d8	10.151	136	912278	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.051	164	575774	20.000	ng/u1	0.00
64) Phenanthrene-d10	16.810	188	1186722	20.000	ng/u1	0.00
79) Chrysene-d12	21.045	240	1034357	20.000	ng/u1	0.00
88) Perylene-d12	23.756	264	1087129	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.022	96	37082	6.303	ng/uL	0.00
4) Pyridine-d5	3.410	84	530708	33.070	ng/u1	0.00
7) Phenol-d5	6.604	99	680592	38.437	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.751	67	429248	36.844	ng/u1	0.00
11) 2-Chlorophenol-d4	6.940	132	546218	36.417	ng/u1	0.00
15) 4-Methylphenol-d8	8.116	113	530839	37.845	ng/u1	0.00
21) Nitrobenzene-d5	8.539	128	278405	38.031	ng/u1	0.00
24) 2-Nitrophenol-d4	9.257	143	315540	39.470	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	9.792	165	555698	36.205	ng/u1	0.00
31) 4-Chloroaniline-d4	10.304	131	748090	37.085	ng/u1	0.00
46) Dimethylphthalate-d6	13.480	166	1609285	36.554	ng/u1	0.00
49) Acenaphthylene-d8	13.739	160	1860124	37.852	ng/u1	0.00
54) 4-Nitrophenol-d4	14.292	143	285462	33.536	ng/u1	0.00
60) Fluorene-d10	15.057	176	1353980	35.953	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.198	200	225213	30.038	ng/u1	0.00
73) Anthracene-d10	16.909	188	2101448	37.354	ng/u1	0.00
81) Pyrene-d10	19.221	212	2321642	39.209	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.574	264	2183538	38.144	ng/u1	0.00

Target Compounds Qvalue

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

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