

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080624\
 Data File : BM047054.D
 Acq On : 07 Aug 2024 12:42
 Operator : RC/JU
 Sample : PB162363BL
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK363

Quant Time: Aug 08 01:13:24 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	7.393	152	236900	20.000	ng/u1	0.00
20) Naphthalene-d8	10.151	136	923161	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.051	164	587325	20.000	ng/u1	0.00
64) Phenanthrene-d10	16.810	188	1191272	20.000	ng/u1	0.00
79) Chrysene-d12	21.039	240	1023847	20.000	ng/u1	-0.01
88) Perylene-d12	23.750	264	1129284	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.016	96	35820	5.990	ng/uL	0.00
4) Pyridine-d5	3.405	84	491565	30.138	ng/u1	0.00
7) Phenol-d5	6.598	99	606668	33.711	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.751	67	386983	32.682	ng/u1	0.00
11) 2-Chlorophenol-d4	6.934	132	485104	31.822	ng/u1	0.00
15) 4-Methylphenol-d8	8.116	113	471645	33.084	ng/u1	0.00
21) Nitrobenzene-d5	8.540	128	242542	32.741	ng/u1	0.00
24) 2-Nitrophenol-d4	9.251	143	275148	34.012	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	9.786	165	484066	31.166	ng/u1	0.00
31) 4-Chloroaniline-d4	10.298	131	654784	32.077	ng/u1	0.00
46) Dimethylphthalate-d6	13.474	166	1432009	31.888	ng/u1	0.00
49) Acenaphthylene-d8	13.739	160	1632894	32.575	ng/u1	0.00
54) 4-Nitrophenol-d4	14.292	143	253468	29.192	ng/u1	0.00
60) Fluorene-d10	15.051	176	1193705	31.073	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.192	200	200932	26.697	ng/u1	0.00
73) Anthracene-d10	16.910	188	1852674	32.806	ng/u1	0.00
81) Pyrene-d10	19.221	212	1995504	34.047	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.568	264	1933378	32.513	ng/u1	0.00

Target Compounds Qvalue

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

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