

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072722\
 Data File : BM035947.D
 Acq On : 27 Jul 2022 12:54
 Operator : CG/JU
 Sample : PB146484BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Jul 27 23:12:25 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM072522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jul 27 03:28:49 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.892	152	314082	20.000	ng/u1	0.00
20) Naphthalene-d8	10.698	136	1320826	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.533	164	792408	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.274	188	1529605	20.000	ng/u1	0.00
79) Chrysene-d12	21.450	240	1301517	20.000	ng/u1	-0.01
88) Perylene-d12	23.838	264	1369007	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	45941	5.504	ng/uL	0.00
4) Pyridine-d5	3.698	84	613914	26.848	ng/u1	0.00
7) Phenol-d5	7.057	99	735986	27.943	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.222	67	459816	26.188	ng/u1	0.00
11) 2-Chlorophenol-d4	7.422	132	649486	29.708	ng/u1	0.00
15) 4-Methylphenol-d8	8.610	113	592427	27.765	ng/u1	-0.01
21) Nitrobenzene-d5	9.063	128	321503	30.718	ng/u1	0.00
24) 2-Nitrophenol-d4	9.786	143	346605	33.991	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.322	165	552719	29.970	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.839	131	870299	27.875	ng/u1	-0.01
46) Dimethylphthalate-d6	13.945	166	1693426	31.373	ng/u1	0.00
49) Acenaphthylene-d8	14.227	160	2161161	31.381	ng/u1	0.00
54) 4-Nitrophenol-d4	14.733	143	295071	26.621	ng/u1	-0.02
60) Fluorene-d10	15.521	176	1455776	29.846	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200	228562	27.799	ng/u1	-0.02
73) Anthracene-d10	17.368	188	2207950	31.928	ng/u1	-0.01
81) Pyrene-d10	19.662	212	2387503	33.099	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.685	264	2249402	32.364	ng/u1	0.00

Target Compounds	Qvalue

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

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