

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072722\
 Data File : BM035949.D
 Acq On : 27 Jul 2022 14:07
 Operator : CG/JU
 Sample : N3709-12
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Jul 27 23:12:43 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	310048	20.000	ng/u1	0.00
20) Naphthalene-d8	10.698	136	1296612	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.527	164	767513	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.268	188	1526070	20.000	ng/u1	-0.01
79) Chrysene-d12	21.444	240	1344348	20.000	ng/u1	-0.02
88) Perylene-d12	23.833	264	1422000	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	16638	2.019	ng/uL	0.00
4) Pyridine-d5	3.687	84	1052	0.047	ng/u1	-0.02
7) Phenol-d5	7.057	99	571611	21.985	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.222	67	435480	25.124	ng/u1	0.00
11) 2-Chlorophenol-d4	7.422	132	500338	23.183	ng/u1	0.00
15) 4-Methylphenol-d8	8.604	113	406757	19.311	ng/u1	-0.02
21) Nitrobenzene-d5	9.057	128	303252	29.515	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.780	143	197377	19.718	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.322	165	412688	22.795	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.839	131	700012	22.840	ng/u1	-0.01
46) Dimethylphthalate-d6	13.945	166	1416802	27.100	ng/u1	0.00
49) Acenaphthylene-d8	14.221	160	2090252	31.336	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.733	143	133771	12.460	ng/u1	-0.02
60) Fluorene-d10	15.521	176	1422815	30.117	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	97053	11.831	ng/u1	-0.02
73) Anthracene-d10	17.368	188	2073310	30.050	ng/u1	-0.01
81) Pyrene-d10	19.656	212	2273559	30.515	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.680	264	2210526	30.620	ng/u1	-0.01

Target Compounds Qvalue

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

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