

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072722\
 Data File : BM035948.D
 Acq On : 27 Jul 2022 13:31
 Operator : CG/JU
 Sample : PB146378BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Jul 27 23:12:34 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	7.893	152	320964	20.000	ng/u1	0.00
20) Naphthalene-d8	10.698	136	1329455	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.527	164	774722	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.268	188	1526920	20.000	ng/u1	-0.01
79) Chrysene-d12	21.445	240	1334724	20.000	ng/u1	-0.02
88) Perylene-d12	23.833	264	1422009	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	54644	6.406	ng/uL	0.00
4) Pyridine-d5	3.699	84	691300	29.584	ng/u1	0.00
7) Phenol-d5	7.057	99	833644	30.972	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.222	67	526127	29.322	ng/u1	0.00
11) 2-Chlorophenol-d4	7.422	132	748227	33.490	ng/u1	0.00
15) 4-Methylphenol-d8	8.604	113	666177	30.551	ng/u1	-0.02
21) Nitrobenzene-d5	9.063	128	362199	34.381	ng/u1	0.00
24) 2-Nitrophenol-d4	9.781	143	386187	37.627	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.322	165	623640	33.596	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.839	131	954311	30.368	ng/u1	-0.01
46) Dimethylphthalate-d6	13.945	166	1857745	35.203	ng/u1	0.00
49) Acenaphthylene-d8	14.227	160	2376697	35.298	ng/u1	0.00
54) 4-Nitrophenol-d4	14.733	143	327169	30.191	ng/u1	-0.02
60) Fluorene-d10	15.521	176	1612745	33.819	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200	255130	31.084	ng/u1	-0.02
73) Anthracene-d10	17.368	188	2425595	35.137	ng/u1	-0.01
81) Pyrene-d10	19.662	212	2683886	36.282	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.680	264	2547324	35.285	ng/u1	-0.01

Target Compounds	Qvalue

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

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