

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072922\
 Data File : BM036014.D
 Acq On : 30 Jul 2022 04:49
 Operator : CG/JU
 Sample : PB146355BL
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK355

Quant Time: Jul 30 05:27:32 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM072522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jul 30 00:20:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	442178	20.000	ng/u1	0.00
20) Naphthalene-d8	10.681	136	1850405	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.510	164	1024743	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.251	188	1861993	20.000	ng/u1	0.00
79) Chrysene-d12	21.427	240	1605280	20.000	ng/u1	0.00
88) Perylene-d12	23.791	264	1658447	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	73736	6.275	ng/uL	0.00
4) Pyridine-d5	3.693	84	972579	30.212	ng/u1	0.00
7) Phenol-d5	7.040	99	1153572	31.110	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.210	67	748797	30.292	ng/u1	0.00
11) 2-Chlorophenol-d4	7.404	132	1014796	32.970	ng/u1	0.00
15) 4-Methylphenol-d8	8.587	113	910163	30.298	ng/u1	0.00
21) Nitrobenzene-d5	9.040	128	486287	33.165	ng/u1	0.00
24) 2-Nitrophenol-d4	9.763	143	496457	34.752	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.298	165	833312	32.252	ng/u1	0.00
31) 4-Chloroaniline-d4	10.822	131	1299637	29.713	ng/u1	0.00
46) Dimethylphthalate-d6	13.927	166	2378778	34.078	ng/u1	0.00
49) Acenaphthylene-d8	14.204	160	3022425	33.936	ng/u1	0.00
54) 4-Nitrophenol-d4	14.710	143	357425	24.935	ng/u1	0.00
60) Fluorene-d10	15.498	176	2040423	32.348	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.621	200	249260	24.904	ng/u1	0.00
73) Anthracene-d10	17.351	188	2891779	34.352	ng/u1	0.00
81) Pyrene-d10	19.639	212	3065757	34.459	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.639	264	2952929	35.072	ng/u1	-0.01

Target Compounds	Qvalue

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

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