

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM083021\
 Data File : BM032012.D
 Acq On : 01 Sep 2021 09:11
 Operator : CG/JU
 Sample : PB138672BL
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK672

Quant Time: Sep 01 10:57:42 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM082821.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Aug 30 09:21:30 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.498	152	92180	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.257	136	398228	20.000	ng/u1	#-0.01
38) Acenaphthene-d10	14.139	164	294928	20.000	ng/u1	-0.01
64) Phenanthrene-d10	16.898	188	648125	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.109	240	672634	20.000	ng/u1	# 0.00
88) Perylene-d12	23.244	264	614010	20.000	ng/u1	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.128	96	12431	6.222	ng/uL	0.00
4) Pyridine-d5	3.516	84	169886	29.945	ng/u1	0.00
7) Phenol-d5	6.698	99	224411	29.358	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.857	67	130045	29.404	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.040	132	185324	31.506	ng/u1	-0.01
15) 4-Methylphenol-d8	8.210	113	190225	29.006	ng/u1	-0.01
21) Nitrobenzene-d5	8.651	128	94760	32.761	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.363	143	109955	32.122	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	9.892	165	208513	30.360	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.416	131	276567	28.548	ng/u1	0.00
46) Dimethylphthalate-d6	13.569	166	716141	32.090	ng/u1	0.00
49) Acenaphthylene-d8	13.827	160	868503	32.862	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.380	143	109197	27.033	ng/u1	0.00
60) Fluorene-d10	15.145	176	625554	31.888	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.292	200	100746	22.622	ng/u1	0.00
73) Anthracene-d10	16.998	188	1017228	33.066	ng/u1	0.00
81) Pyrene-d10	19.303	212	1167616	37.061	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.109	264	1056972	32.754	ng/u1	0.00

Target Compounds	Qvalue

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

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