

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052021\
 Data File : BM030151.D
 Acq On : 22 May 2021 15:39
 Operator : CG/JU
 Sample : M2455-01
 Misc :
 ALS Vial : 67 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0BA1

Quant Time: May 24 02:30:37 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050721.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon May 24 02:23:05 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.50	152	140154	20.00	ng/ul	-0.01
20) Naphthalene-d8	10.27	136	561401	20.00	ng/ul	-0.01
38) Acenaphthene-d10	14.16	164	387120	20.00	ng/ul	0.00
64) Phenanthrene-d10	16.91	188	789110	20.00	ng/ul	0.00
79) Chrysene-d12	21.10	240	817322	20.00	ng/ul	0.00
88) Perylene-d12	23.24	264	768822	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL
4) Pyridine-d5	0.00	84	0d	0.00	ng/ul
7) Phenol-d5	0.00	99	0	0.00	ng/ul
9) Bis-(2-Chloroethyl)ether-d	0.00	67	0	0.00	ng/ul
11) 2-Chlorophenol-d4	0.00	132	0	0.00	ng/ul
15) 4-Methylphenol-d8	0.00	113	0	0.00	ng/ul
21) Nitrobenzene-d5	0.00	128	0	0.00	ng/ul
24) 2-Nitrophenol-d4	0.00	143	0	0.00	ng/ul
28) 2,4-Dichlorophenol-d3	0.00	165	0	0.00	ng/ul
31) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul
46) Dimethylphthalate-d6	0.00	166	0	0.00	ng/ul
49) Acenaphthylene-d8	0.00	160	0	0.00	ng/ul
54) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul
60) Fluorene-d10	0.00	176	0	0.00	ng/ul
65) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul
73) Anthracene-d10	0.00	188	0d	0.00	ng/ul
81) Pyrene-d10	0.00	212	0d	0.00	ng/ul
92) Benzo(a)pyrene-d12	0.00	264	0d	0.00	ng/ul

Target Compounds				Ovalue
72) Phenanthrene	16.95	178	50304	1.081 ng/ul 98
80) Fluoranthene	18.97	202	80294	1.574 ng/ul# 96
87) Chrysene	21.14	228	54701	1.030 ng/ul 97

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

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