

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM060120\
 Data File : BM026223.D
 Acq On : 02 Jun 2020 12:07
 Operator : JU/CG
 Sample : L2829-10DL 10X
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C5CE3DL

Quant Time: Jun 02 12:45:37 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jun 02 03:00:44 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.47	152	154879	20.00	ng/ul	0.00
18) Naphthalene-d8	10.23	136	609766	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.12	164	364585	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.87	188	726002	20.00	ng/ul	0.00
78) Chrysene-d12	21.07	240	752747	20.00	ng/ul	0.00
86) Perylene-d12	23.18	264	797879	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
5) Phenol-d5	6.68	99	9790	0.70	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.82	67	25143	2.59	ng/ul	0.00
9) 2-Chlorophenol-d4	7.01	132	23010	2.09	ng/ul	0.00
13) 4-Methylphenol-d8	8.19	113	15874	1.50	ng/ul	0.00
19) Nitrobenzene-d5	8.62	128	13067	2.63	ng/ul	0.00
22) 2-Nitrophenol-d4	9.33	143	13503	2.75	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.87	165	25273	2.56	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul	
44) Dimethylphthalate-d6	13.54	166	97269	3.33	ng/ul	0.00
47) Acenaphthylene-d8	13.80	160	110657	3.19	ng/ul	0.00
52) 4-Nitrophenol-d4	0.00	143	0d	0.00	ng/ul	
58) Fluorene-d10	15.12	176	82947	3.26	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.26	200	11372	2.69	ng/ul	0.00
71) Anthracene-d10	16.97	188	123681	3.46	ng/ul	0.00
79) Pyrene-d10	19.26	212	133252	3.45	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.04	264	115174	2.71	ng/ul	0.00

Target Compounds						Qvalue
56) 2,3,4,6-Tetrachlorophenol	14.76	232	475495	67.40	ng/ul#	99
69) Pentachlorophenol	16.56	266	7219178	1378.93	ng/ul	98

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

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