

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
 Data File : BM026527.D
 Acq On : 24 Jun 2020 09:33
 Operator : JU/CG
 Sample : L3069-11DL 2X
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 CB7H0DL

Quant Time: Jun 24 10:32:01 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jun 23 10:25:31 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.37	152	103857	20.00	ng/ul	0.00
18) Naphthalene-d8	10.13	136	473591	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.03	164	311839	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.79	188	704519	20.00	ng/ul	0.00
78) Chrysene-d12	21.01	240	761230	20.00	ng/ul	0.00
86) Perylene-d12	23.09	264	709536	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	6.59	99	21252	2.22	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.73	67	59518	9.83	ng/ul	0.00
9) 2-Chlorophenol-d4	6.92	132	59511	7.63	ng/ul	0.00
13) 4-Methylphenol-d8	8.10	113	42775	5.68	ng/ul	0.00
19) Nitrobenzene-d5	8.52	128	33357	7.93	ng/ul	0.00
22) 2-Nitrophenol-d4	9.24	143	38118	8.23	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.78	165	71016	8.33	ng/ul	0.00
29) 4-Chloroaniline-d4	10.29	131	97222	10.94	ng/ul	0.00
44) Dimethylphthalate-d6	13.46	166	241792	9.22	ng/ul	0.00
47) Acenaphthylene-d8	13.72	160	243365	7.74	ng/ul	0.00
52) 4-Nitrophenol-d4	0.00	143	0d	0.00	ng/ul	
58) Fluorene-d10	15.04	176	182307	8.10	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.19	200	20200	4.38	ng/ul	0.00
71) Anthracene-d10	16.89	188	286930	8.00	ng/ul	0.00
79) Pyrene-d10	19.20	212	336168	8.05	ng/ul	0.00
90) Benzo(a)pyrene-d12	22.96	264	216453	5.52	ng/ul	0.00

Target Compounds					Ovalue	
2) 1,4-Dioxane	2.99	88	84040	25.596	ng/uL	95
45) Dimethylphthalate	13.51	163	92630	3.402	ng/ul	99

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

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