

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062520\
 Data File : BM026565.D
 Acq On : 25 Jun 2020 19:38
 Operator : JU/CG
 Sample : L3062-16
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ET149

Manual Integrations
 APPROVED

mohammad
 6/26/2020 12:20:11 PM

Quant Time: Jun 25 20:11:33 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jun 25 10:53:13 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.37	152	92765	20.00	ng/ul	0.00
18) Naphthalene-d8	10.13	136	445592	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.03	164	309635	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.79	188	723386	20.00	ng/ul	0.00
78) Chrysene-d12	21.00	240	808206	20.00	ng/ul	0.00
86) Perylene-d12	23.08	264	804601	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.97	96	6288	2.38	ng/uL	0.00
5) Phenol-d5	6.58	99	90531	10.59	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.73	67	70916	13.12	ng/ul	0.00
9) 2-Chlorophenol-d4	6.91	132	74362	10.67	ng/ul	0.00
13) 4-Methylphenol-d8	8.10	113	78294	11.64	ng/ul	0.00
19) Nitrobenzene-d5	8.52	128	33560m	8.48	ng/ul	0.00
22) 2-Nitrophenol-d4	9.23	143	27658	6.35	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.77	165	70147	8.75	ng/ul	0.00
29) 4-Chloroaniline-d4	10.29	131	95793	11.45	ng/ul	0.00
44) Dimethylphthalate-d6	13.46	166	276345	10.61	ng/ul	0.00
47) Acenaphthylene-d8	13.72	160	232520	7.45	ng/ul	0.00
52) 4-Nitrophenol-d4	14.32	143	6207m	1.44	ng/ul	0.05
58) Fluorene-d10	15.03	176	153905	6.89	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.19	200	6163m	1.30	ng/ul	0.01
71) Anthracene-d10	16.89	188	211911	5.75	ng/ul	0.00
79) Pyrene-d10	19.19	212	215758	4.87	ng/ul	0.00
90) Benzo(a)pyrene-d12	22.95	264	165754	3.73	ng/ul	0.00

Target Compounds

					Ovalue
16) 4-Methylphenol	8.15	108	19724	2.678	ng/ul 94
45) Dimethylphthalate	13.50	163	201750	7.462	ng/ul 99

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

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