

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062117\
 Data File : BM010635.D
 Acq On : 21 Jun 2017 22:28
 Operator : SJ/JU
 Sample : I3627-14
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F5E25

Manual Integrations
 APPROVED

mohammad
 6/22/2017 5:32:59 PM

Quant Time: Jun 22 01:48:52 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM062017.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jun 22 01:38:18 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.46	152	76433	20.00	ng/ul	0.00
18) Naphthalene-d8	10.23	136	346815	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.12	164	241238	20.00	ng/ul	0.00
61) Phenanthrene-d10	16.87	188	607095	20.00	ng/ul	0.00
75) Chrysene-d12	21.09	240	729914	20.00	ng/ul	0.00
83) Perylene-d12	23.23	264	655950	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.07	96	5275	3.95	ng/uL	0.00
5) Phenol-d5	6.65	99	119122	16.36	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.82	67	72273	17.39	ng/ul	0.00
9) 2-Chlorophenol-d4	7.00	132	94161	18.43	ng/ul	0.00
13) 4-Methylphenol-d8	8.17	113	103959	17.22	ng/ul	0.00
19) Nitrobenzene-d5	8.60	128	51557	20.84	ng/ul	0.00
22) 2-Nitrophenol-d4	9.32	143	60314	21.16	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.85	165	117508	19.58	ng/ul	0.00
29) 4-Chloroaniline-d4	10.37	131	140352	22.19	ng/ul	0.00
43) Dimethylphthalate-d6	13.54	166	479788	22.66	ng/ul	0.00
46) Acenaphthylene-d8	13.80	160	528148	21.44	ng/ul	0.00
51) 4-Nitrophenol-d4	14.32	143	51921	13.93	ng/ul	0.00
57) Fluorene-d10	15.12	176	401340	21.92	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.24	200	63578	17.05	ng/ul	0.00
70) Anthracene-d10	16.97	188	651180	22.21	ng/ul	0.00
76) Pyrene-d10	19.28	212	783516	23.11	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.09	264	716778m	22.61	ng/ul	0.00
Target Compounds						
6) Phenol	6.67	94	10896	1.453	ng/ul#	86
44) Dimethylphthalate	13.59	163	225314	10.854	ng/ul	97
74) Fluoranthene	18.94	202	119562	2.874	ng/ul#	95
77) Pyrene	19.30	202	144699	3.432	ng/ul#	93
80) Benzo(a)anthracene	21.07	228	95719	2.252	ng/ul	96
82) Chrysene	21.12	228	92453m	2.440	ng/ul	
85) Benzo(b)fluoranthene	22.59	252	147965	3.475	ng/ul#	97
86) Benzo(k)fluoranthene	22.63	252	54079m	1.340	ng/ul	
88) Benzo(a)pyrene	23.13	252	71280	1.818	ng/ul	97
89) Indeno(1,2,3-cd)pyrene	25.34	276	40761	1.003	ng/ul#	94

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

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