

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080719\
 Data File : BG042040.D
 Acq On : 7 Aug 2019 20:47
 Operator : HP/JU
 Sample : K4029-19
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BENT8

Manual Integrations
 APPROVED

mohammad
 8/9/2019 4:30:19 PM

Quant Time: Aug 08 02:55:24 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080319MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Aug 08 01:36:24 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.03	152	76813	20.00	ng/ul	0.00
18) Naphthalene-d8	10.85	136	327971	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.69	164	233783	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.44	188	491523	20.00	ng/ul	0.00
77) Chrysene-d12	21.72	240	483746	20.00	ng/ul	0.00
85) Perylene-d12	24.99	264	570949	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.41	96	6964	3.97	ng/uL	0.00
5) Phenol-d5	7.17	99	151481	25.13	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.34	67	86747	25.73	ng/ul	0.00
9) 2-Chlorophenol-d4	7.55	132	141087	27.38	ng/ul	0.00
13) 4-Methylphenol-d8	8.73	113	123470	24.40	ng/ul	0.00
19) Nitrobenzene-d5	9.19	128	73508	29.77	ng/ul	0.00
22) 2-Nitrophenol-d4	9.92	143	92490	32.27	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.46	165	164338	28.40	ng/ul	0.00
29) 4-Chloroaniline-d4	10.98	131	165579	25.79	ng/ul	0.00
43) Dimethylphthalate-d6	14.08	166	571843	31.54	ng/ul	0.00
46) Acenaphthylene-d8	14.38	160	606812	29.30	ng/ul	0.00
51) 4-Nitrophenol-d4	14.87	143	65279	21.44	ng/ul	0.00
57) Fluorene-d10	15.68	176	488436	30.28	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.79	200	80691	25.59	ng/ul	0.00
70) Anthracene-d10	17.54	188	744703	31.58	ng/ul	0.00
78) Pyrene-d10	19.82	212	819890	30.37	ng/ul	0.00
89) Benzo(a)pyrene-d12	24.77	264	932554	31.27	ng/ul	0.00

Target Compounds

					Ovalue	
6) Phenol	7.20	94	15297	2.458	ng/ul#	93
44) Dimethylphthalate	14.13	163	218031	12.111	ng/ul	99
69) Phenanthrene	17.48	178	38865	1.444	ng/ul	98
76) Fluoranthene	19.49	202	109069	3.508	ng/ul	96
79) Pyrene	19.85	202	103508	3.181	ng/ul	97
82) Benzo(a)anthracene	21.70	228	68360	2.038	ng/ul	97
84) Chrysene	21.77	228	70473	2.251	ng/ul	97
87) Benzo(b)fluoranthene	23.95	252	86989	2.428	ng/ul#	98
88) Benzo(k)fluoranthene	24.01	252	37408m	1.086	ng/ul	
90) Benzo(a)pyrene	24.83	252	63367	1.856	ng/ul	97
91) Indeno(1,2,3-cd)pyrene	28.70	276	40502	1.017	ng/ul	97
93) Benzo(g,h,i)perylene	29.87	276	33491m	1.007	ng/ul	

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

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