

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080315\
 Data File : BM002203.D
 Acq On : 03 Aug 2015 18:35
 Operator : TP/UM
 Sample : G3095-23
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C01P0

Manual Integrations
 APPROVED

MMDadoda
 8/6/2015 4:15:54 PM

Quant Time: Aug 04 04:36:07 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 04 02:31:26 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.53	152	98606	20.00	ng/ul	0.00
18) Naphthalene-d8	10.32	136	385486	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.21	164	204382	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.96	188	428114	20.00	ng/ul	0.00
78) Chrysene-d12	21.17	240	510848	20.00	ng/ul	0.00
86) Perylene-d12	23.36	264	551461	20.00	ng/ul	0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.98	96	2768	1.45	ng/uL	0.00
5) Phenol-d5	6.73	99	67038	9.51	ng/ul	0.01
7) Bis-(2-Chloroethyl)ether-d	6.87	67	36589	9.39	ng/ul	0.00
9) 2-Chlorophenol-d4	7.07	132	59446	9.95	ng/ul	0.00
13) 4-Methylphenol-d8	8.26	113	51417	9.17	ng/ul	0.01
19) Nitrobenzene-d5	8.69	128	27649	10.09	ng/ul	0.00
22) 2-Nitrophenol-d4	9.42	143	31945	10.49	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.96	165	57084	9.73	ng/ul	0.00
29) 4-Chloroaniline-d4	10.47	131	42472	6.59	ng/ul	0.00
44) Dimethylphthalate-d6	13.62	166	152353	10.18	ng/ul	0.00
47) Acenaphthylene-d8	13.90	160	184293	9.85	ng/ul	0.00
52) 4-Nitrophenol-d4	14.46	143	15097	6.04	ng/ul	0.02
58) Fluorene-d10	15.21	176	125809	9.55	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.36	200	3709	1.42	ng/ul	0.01
71) Anthracene-d10	17.06	188	177369	9.28	ng/ul	0.00
79) Pyrene-d10	19.37	212	193497	9.00	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.22	264	228229	8.54	ng/ul	0.01

Target Compounds

						Qvalue
45) Dimethylphthalate	13.67	163	58499	3.91	ng/ul	99
70) Phenanthrene	17.00	178	114689	5.17	ng/ul	99
72) Anthracene	17.09	178	35913	1.58	ng/ul	97
77) Fluoranthene	19.03	202	246803	9.64	ng/ul	100
80) Pyrene	19.40	202	254987	9.17	ng/ul	99
83) Benzo(a)anthracene	21.16	228	118289	4.18	ng/ul	94
84) Bis(2-ethylhexyl)phthalate	21.08	149	170033	10.46	ng/ul	98
85) Chrysene	21.21	228	138961	5.24	ng/ul	97
88) Benzo(b)fluoranthene	22.71	252	180107	5.81	ng/ul#	92
89) Benzo(k)fluoranthene	22.74	252	65895m	2.20	ng/ul	
91) Benzo(a)pyrene	23.27	252	111043	3.71	ng/ul#	94
92) Indeno(1,2,3-cd)pyrene	25.56	276	82127	2.38	ng/ul	94
94) Benzo(g,h,i)perylene	26.24	276	81115	2.81	ng/ul	98

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

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