

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112119\
 Data File : BM023833.D
 Acq On : 21 Nov 2019 17:31
 Operator : JU
 Sample : K5851-10
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AA8

Manual Integrations
 APPROVED

mohammad
 11/22/2019 1:22:42 PM

Quant Time: Nov 22 05:57:03 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111119MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 21 16:09:32 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.51	152	328624	20.00	ng/ul	-0.01
18) Naphthalene-d8	10.29	136	1352211	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.17	164	830256	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.92	188	1783854	20.00	ng/ul	0.00
78) Chrysene-d12	21.13	240	1745304	20.00	ng/ul	0.00
86) Perylene-d12	23.29	264	2116623	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.06	96	14821	1.86	ng/uL	0.00
5) Phenol-d5	6.70	99	406889	13.58	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.86	67	233865	13.59	ng/ul	0.00
9) 2-Chlorophenol-d4	7.05	132	324892	14.97	ng/ul	0.00
13) 4-Methylphenol-d8	8.23	113	315748	13.21	ng/ul	-0.01
19) Nitrobenzene-d5	8.66	128	153509	15.86	ng/ul	0.00
22) 2-Nitrophenol-d4	9.38	143	167376	16.45	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.92	165	322036	13.88	ng/ul	0.00
29) 4-Chloroaniline-d4	10.43	131	318755	12.51	ng/ul	0.00
44) Dimethylphthalate-d6	13.59	166	967511	14.94	ng/ul	0.00
47) Acenaphthylene-d8	13.86	160	1161937	15.73	ng/ul	0.00
52) 4-Nitrophenol-d4	14.41	143	92669	8.38	ng/ul	0.00
58) Fluorene-d10	15.17	176	887697	15.53	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.31	200	65627	6.36	ng/ul	0.00
71) Anthracene-d10	17.02	188	1431087	16.91	ng/ul	0.00
79) Pyrene-d10	19.33	212	1558628	17.13	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.15	264	1885831	16.93	ng/ul	0.00

Target Compounds

					Ovalue	
28) Naphthalene	10.33	128	118856	1.672	ng/ul	97
45) Dimethylphthalate	13.64	163	456915	7.201	ng/ul	100
48) Acenaphthylene	13.89	152	548206	7.370	ng/ul	100
50) Acenaphthene	14.23	153	115231	2.183	ng/ul	100
54) Dibenzofuran	14.57	168	334257	4.277	ng/ul	98
59) Fluorene	15.23	166	746663	11.778	ng/ul	100
70) Phenanthrene	16.97	178	7843887	79.830	ng/ul	99
72) Anthracene	17.06	178	1878351	19.012	ng/ul	100
75) Carbazole	17.33	167	445512	5.409	ng/ul	99
77) Fluoranthene	19.00	202	11468943m	108.947	ng/ul	
80) Pyrene	19.36	202	8863842	78.571	ng/ul	99
83) Benzo(a)anthracene	21.12	228	4450741	38.661	ng/ul	97
84) Bis(2-ethylhexyl)phthalate	21.07	149	71746	1.330	ng/ul#	97
85) Chrysene	21.17	228	4120410	38.018	ng/ul	97
88) Benzo(b)fluoranthene	22.65	252	5491350	40.698	ng/ul	98
89) Benzo(k)fluoranthene	22.69	252	1891865m	14.554	ng/ul	
91) Benzo(a)pyrene	23.19	252	3943769	31.928	ng/ul	97
92) Indeno(1,2,3-cd)pyrene	25.43	276	2712403	20.536	ng/ul	98
93) Dibenzo(a,h)anthracene	25.43	278	713861	6.404	ng/ul#	83
94) Benzo(g,h,i)perylene	26.07	276	2567230	23.986	ng/ul	97

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

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