

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120120\
 Data File : BG047529.D
 Acq On : 2 Dec 2020 13:21
 Operator : CG/JU
 Sample : L4865-05
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0B77

Manual Integrations
 APPROVED

mohammad
 12/3/2020 3:34:57 PM

Quant Time: Dec 02 14:18:04 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 23 13:14:53 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.26	152	28766	20.00	ng/ul	-0.01
18) Naphthalene-d8	11.09	136	114675	20.00	ng/ul	-0.01
36) Acenaphthene-d10	14.87	164	90563	20.00	ng/ul	-0.01
62) Phenanthrene-d10	17.61	188	222741	20.00	ng/ul	0.00
78) Chrysene-d12	21.89	240	222232	20.00	ng/ul	0.00
86) Perylene-d12	25.29	264	236213	20.00	ng/ul	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.61	96	1656	2.09	ng/uL	0.00
5) Phenol-d5	7.40	99	38145	13.22	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.58	67	22575	13.72	ng/ul	0.00
9) 2-Chlorophenol-d4	7.79	132	31015	16.03	ng/ul	-0.01
13) 4-Methylphenol-d8	8.95	113	25730	11.61	ng/ul	-0.01
19) Nitrobenzene-d5	9.43	128	16236	17.14	ng/ul	0.00
22) 2-Nitrophenol-d4	10.16	143	17700	17.13	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.70	165	38103	16.10	ng/ul	-0.01
29) 4-Chloroaniline-d4	11.22	131	32360	13.52	ng/ul	0.00
44) Dimethylphthalate-d6	14.27	166	134106	17.25	ng/ul	-0.01
47) Acenaphthylene-d8	14.57	160	157824	17.57	ng/ul	0.00
52) 4-Nitrophenol-d4	15.03	143	17955	15.14	ng/ul	-0.01
58) Fluorene-d10	15.86	176	129900	18.24	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.97	200	24376	19.23	ng/ul	0.00
71) Anthracene-d10	17.71	188	203976	18.46	ng/ul	-0.01
79) Pyrene-d10	19.97	212	257951	19.95	ng/ul	0.00
90) Benzo(a)pyrene-d12	25.06	264	264672	19.47	ng/ul	-0.02

Target Compounds				Ovalue		
6) Phenol	7.42	94	4978	1.801	ng/ul#	75
45) Dimethylphthalate	14.32	163	29188	3.681	ng/ul#	96
70) Phenanthrene	17.65	178	162221	13.486	ng/ul#	95
72) Anthracene	17.75	178	18258m	1.507	ng/ul	
75) Carbazole	18.01	167	16504	1.689	ng/ul	96
77) Fluoranthene	19.64	202	271047	17.843	ng/ul#	98
80) Pyrene	20.00	202	200563	12.741	ng/ul#	97
83) Benzo(a)anthracene	21.87	228	91181	5.766	ng/ul	99
85) Chrysene	21.94	228	100917	6.707	ng/ul	98
88) Benzo(b)fluoranthene	24.21	252	110939	6.652	ng/ul	95
89) Benzo(k)fluoranthene	24.27	252	50215	3.055	ng/ul	92
91) Benzo(a)pyrene	25.13	252	70403m	4.735	ng/ul	
92) Indeno(1,2,3-cd)pyrene	29.18	276	49059m	2.713	ng/ul	
94) Benzo(g,h,i)perylene	30.40	276	39909m	2.674	ng/ul	

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

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