

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082319\
 Data File : BM022250.D
 Acq On : 23 Aug 2019 11:28
 Operator : HP/JU
 Sample : K4183-20
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AG5

Manual Integrations
 APPROVED

mohammad
 8/26/2019 8:29:24 AM

Quant Time: Aug 23 13:09:24 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM082219MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Aug 22 19:41:45 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	35538	20.00	ng/ul	0.00
18) Naphthalene-d8	10.33	136	156110	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.22	164	108491	20.00	ng/ul	0.00
61) Phenanthrene-d10	16.97	188	250120	20.00	ng/ul	0.00
77) Chrysene-d12	21.19	240	289834	20.00	ng/ul	0.00
85) Perylene-d12	23.38	264	259096	20.00	ng/ul	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.15	96	1429	1.80	ng/uL	0.01
5) Phenol-d5	6.75	99	32389	10.56	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.92	67	20852	11.69	ng/ul	0.00
9) 2-Chlorophenol-d4	7.10	132	28439	11.48	ng/ul	0.00
13) 4-Methylphenol-d8	8.28	113	25126	9.45	ng/ul	0.00
19) Nitrobenzene-d5	8.73	128	13838	11.89	ng/ul	0.00
22) 2-Nitrophenol-d4	9.45	143	15671	11.57	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.98	165	31617	11.46	ng/ul	0.00
29) 4-Chloroaniline-d4	10.51	131	29098	8.67	ng/ul	0.01
43) Dimethylphthalate-d6	13.64	166	126646	13.26	ng/ul	0.00
46) Acenaphthylene-d8	13.91	160	131261	12.75	ng/ul	0.00
51) 4-Nitrophenol-d4	14.49	143	9539	6.74	ng/ul	0.02
57) Fluorene-d10	15.22	176	106658	13.46	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.39	200	15356	7.96	ng/ul	0.00
70) Anthracene-d10	17.08	188	178913	13.51	ng/ul	0.00
78) Pyrene-d10	19.39	212	240932	15.99	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.24	264	203344	14.29	ng/ul	0.00

Target Compounds				Ovalue		
6) Phenol	6.77	94	4048	1.233	ng/ul#	83
44) Dimethylphthalate	13.69	163	31374	3.227	ng/ul	98
69) Phenanthrene	17.02	178	102985	6.882	ng/ul	98
71) Anthracene	17.12	178	30568	1.963	ng/ul	99
76) Fluoranthene	19.05	202	167288	7.419	ng/ul	99
79) Pyrene	19.42	202	139153	7.418	ng/ul	97
82) Benzo(a)anthracene	21.17	228	83477	3.827	ng/ul	96
84) Chrysene	21.22	228	60590	2.971	ng/ul	97
87) Benzo(b)fluoranthene	22.73	252	78257	4.342	ng/ul#	95
88) Benzo(k)fluoranthene	22.76	252	26103	1.493	ng/ul#	97
90) Benzo(a)pyrene	23.29	252	55514	3.231	ng/ul#	92
91) Indeno(1,2,3-cd)pyrene	25.57	276	31114m	1.560	ng/ul	
93) Benzo(g,h,i)perylene	26.25	276	27247m	1.790	ng/ul	

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

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082319\
Data File : BM022250.D
Acq On : 23 Aug 2019 11:28
Operator : HP/JU
Sample : K4183-20
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleID :
C0AG5

Quant Time: Aug 23 13:09:24 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM082219MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Aug 22 19:41:45 2019
Response via : Initial Calibration

Manual Integrations
APPROVED

mohammad
8/26/2019 8:29:24 AM

