Quantitation Report (QT/LSC Reviewed)

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM101719\
Data File : BM023241.D
Acq On : 18 Oct 2019 04:41
Operator : JU
Sample : K5262-18DL 5X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFB79DL

Manual Integrations APPROVED
mohammad
10/20/2019 8:06:44 PM

Quant Time: Oct 18 07:55:26 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM101419MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Oct 18 07:48:41 2019
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	441740	20.00	ng/ul	0.00
18) Naphthalene-d8	10.44	136	1739390	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.30	164	1105542	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.05	188	2349728	20.00	ng/ul	0.00
78) Chrysene-d12	21.24	240	2521298	20.00	ng/ul	0.00
86) Perylene-d12	23.45	264	3011008	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.18	96	3566	0.36	ng/uL	0.00
5) Phenol-d5	6.83	99	50999	1.33	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.00	67	29574	1.36	ng/ul	0.00
9) 2-Chlorophenol-d4	7.19	132	43241	1.47	ng/ul	0.00
13) 4-Methylphenol-d8	8.36	113	40153	1.30	ng/ul	0.00
19) Nitrobenzene-d5	8.80	128	20441	1.74	ng/ul	0.00
22) 2-Nitrophenol-d4	9.52	143	19901	1.78	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.06	165	42948	1.51	ng/ul	0.00
29) 4-Chloroaniline-d4	10.57	131	31885	0.92	ng/ul	0.00
44) Dimethylphthalate-d6	13.72	166	144227	1.70	ng/ul	0.00
47) Acenaphthylene-d8	14.00	160	163671	1.65	ng/ul	0.00
52) 4-Nitrophenol-d4	14.50	143	12418	0.91	ng/ul	0.00
58) Fluorene-d10	15.30	176	130367	1.81	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.42	200	7626	0.82	ng/ul	0.00
71) Anthracene-d10	17.15	188	212336	1.87	ng/ul	0.00
79) Pyrene-d10	19.44	212	241833	1.98	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.31	264	285602	1.87	ng/ul	0.00
Target Compounds						Qvalue
48) Acenaphthylene	14.02	152	139415	1.372	ng/ul	98
50) Acenaphthene	14.37	153	105716	1.494	ng/ul	98
54) Dibenzofuran	14.70	168	131757	1.312	ng/ul	98
59) Fluorene	15.36	166	140293	1.703	ng/ul	99
70) Phenanthrene	17.09	178	2697046	20.453	ng/ul	100
72) Anthracene	17.18	178	578420	4.245	ng/ul	100
75) Carbazole	17.45	167	266230	2.187	ng/ul	98
77) Fluoranthene	19.12	202	4581837	29.215	ng/ul	98
80) Pyrene	19.47	202	3903644	24.715	ng/ul	100
83) Benzo(a)anthracene	21.23	228	2136121	12.504	ng/ul	98
85) Chrysene	21.27	228	2052768	12.941	ng/ul	97
88) Benzo(b)fluoranthene	22.79	252	3186935	17.457	ng/ul	98
89) Benzo(k)fluoranthene	22.83	252	1012958m	5.779	ng/ul	
91) Benzo(a)pyrene	23.36	252	2252774	12.862	ng/ul	97
92) Indeno(1,2,3-cd)pyrene	25.67	276	1763152	8.462	ng/ul	97
93) Dibenzo(a,h)anthracene	25.67	278	447232	2.567	ng/ul#	91
94) Benzo(g,h,i)perylene	26.34	276	1224340	7.137	ng/ul	99

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

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