

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082019\
 Data File : BG042368.D
 Acq On : 20 Aug 2019 21:52
 Operator : HP/JU
 Sample : K4386-20
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 T6-B-(0-2)

Manual Integrations
 APPROVED

mohammad
 8/22/2019 12:49:20 PM

Quant Time: Aug 21 05:34:33 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG081919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 19 13:40:37 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.01	152	71807	20.00	ng	0.00
21) Naphthalene-d8	10.83	136	275583	20.00	ng	0.00
39) Acenaphthene-d10	14.66	164	197811	20.00	ng	0.00
64) Phenanthrene-d10	17.41	188	391975	20.00	ng	0.00
76) Chrysene-d12	21.69	240	407826	20.00	ng	0.00
87) Perylene-d12	24.94	264	483260	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	369481	104.25	ng	0.00
7) Phenol-d6	7.17	99	518186	107.66	ng	0.00
23) Nitrobenzene-d5	9.17	82	293782	74.98	ng	0.00
42) 2,4,6-Tribromophenol	16.15	330	274771	101.99	ng	0.00
45) 2-Fluorobiphenyl	13.28	172	719783	61.57	ng	0.00
79) Terphenyl-d14	20.03	244	1041893	55.96	ng	0.00
Target Compounds						
15) Benzyl Alcohol	8.25	79	28024	8.466	ng	96
50) Dimethylphthalate	14.11	163	47486	3.378	ng	98
71) Phenanthrene	17.46	178	235513	12.064	ng	98
72) Anthracene	17.55	178	53448	2.745	ng	97
75) Fluoranthene	19.47	202	418989	17.798	ng	99
78) Pyrene	19.83	202	381819	15.215	ng	98
81) Benzo(a)anthracene	21.67	228	220049	8.786	ng	97
83) Chrysene	21.74	228	197498	8.243	ng	97
84) Bis(2-ethylhexyl)phthalate	21.58	149	41189	2.970	ng	98
86) Indeno(1,2,3-cd)pyrene	28.63	276	145238	4.540	ng	97
88) Benzo(b)fluoranthene	23.90	252	280603	10.564	ng	99
89) Benzo(k)fluoranthene	23.96	252	110125m	4.241	ng	
90) Benzo(a)pyrene	24.78	252	205286	8.065	ng	95
92) Benzo(a,h,i)perylene	29.79	276	145403	5.746	ng	97

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

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