

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112019\
 Data File : BF117889.D
 Acq On : 20 Nov 2019 13:31
 Operator : JU
 Sample : K5908-03
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B-8-2-(6-12)

Manual Integrations
 APPROVED

mohammad
 11/21/2019 1:05:06 PM

Quant Time: Nov 21 03:22:05 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 13 18:36:07 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.92	152	154823	20.00	ng	-0.01
21) Naphthalene-d8	8.20	136	584779	20.00	ng	-0.01
39) Acenaphthene-d10	9.96	164	333739	20.00	ng	-0.01
64) Phenanthrene-d10	11.46	188	635891	20.00	ng	-0.01
76) Chrysene-d12	14.10	240	374981	20.00	ng	-0.01
87) Perylene-d12	15.62	264	423566	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	974864	111.46	ng	0.00
7) Phenol-d6	6.54	99	1299999	123.23	ng	0.00
23) Nitrobenzene-d5	7.47	82	793027	80.62	ng	-0.02
42) 2,4,6-Tribromophenol	10.76	330	394574	111.67	ng	-0.01
45) 2-Fluorobiphenyl	9.28	172	1478987	79.50	ng	-0.01
79) Terphenyl-d14	13.05	244	1315605	64.12	ng	0.00
Target Compounds						
10) Phenol	6.55	94	72597	6.622	ng	86
50) Dimethylphthalate	9.67	163	257442	11.000	ng	97
58) Fluorene	10.52	166	48964	2.399	ng	97
71) Phenanthrene	11.48	178	764399	23.910	ng	97
72) Anthracene	11.53	178	143333	4.522	ng	96
73) Carbazole	11.69	167	72175	2.606	ng	94
75) Fluoranthene	12.67	202	702106	19.923	ng	99
78) Pyrene	12.90	202	490794	16.196	ng	98
81) Benzo(a)anthracene	14.09	228	246907	9.964	ng	99
83) Chrysene	14.13	228	222156	9.252	ng	97
86) Indeno(1,2,3-cd)pyrene	17.14	276	89954	4.402	ng	96
88) Benzo(b)fluoranthene	15.16	252	280858m	10.206	ng	
89) Benzo(k)fluoranthene	15.19	252	73346m	2.829	ng	
90) Benzo(a)pyrene	15.54	252	172874	7.144	ng	98
92) Benzo(a,h,i)perylene	17.60	276	76555	3.690	ng	99

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

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