

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112519\
 Data File : BF117987.D
 Acq On : 26 Nov 2019 1:14
 Operator : JU
 Sample : K5970-03
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-28-S

Manual Integrations
 APPROVED

mohammad
 11/26/2019 12:49:23 PM

Quant Time: Nov 26 05:45:56 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 26 05:38:20 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	193704	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	685793	20.00	ng	0.00
39) Acenaphthene-d10	9.94	164	393612	20.00	ng	0.00
64) Phenanthrene-d10	11.43	188	545606	20.00	ng	0.00
76) Chrysene-d12	14.08	240	284712	20.00	ng	0.00
87) Perylene-d12	15.58	264	356561	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.51	112	1061922	97.04	ng	0.01
7) Phenol-d6	6.52	99	1513553	114.67	ng	0.00
23) Nitrobenzene-d5	7.46	82	833142	72.22	ng	0.00
42) 2,4,6-Tribromophenol	10.73	330	430904	103.41	ng	0.00
45) 2-Fluorobiphenyl	9.26	172	1720466	78.42	ng	0.00
79) Terphenyl-d14	13.02	244	1363714	87.54	ng	0.00
Target Compounds						
49) Acenaphthylene	9.80	152	132476	3.782	ng	97
50) Dimethylphthalate	9.65	163	266209	9.645	ng	99
71) Phenanthrene	11.46	178	436322	15.906	ng	96
72) Anthracene	11.50	178	106872	3.929	ng	96
75) Fluoranthene	12.65	202	1008290	39.433	ng	99
78) Pyrene	12.88	202	983223	42.732	ng	99
81) Benzo(a)anthracene	14.07	228	528641	28.098	ng	94
83) Chrysene	14.11	228	480929	26.380	ng	100
86) Indeno(1,2,3-cd)pyrene	17.10	276	225770	14.550	ng	95
88) Benzo(b)fluoranthene	15.14	252	625612m	27.006	ng	
89) Benzo(k)fluoranthene	15.16	252	211138m	9.673	ng	
90) Benzo(a)pyrene	15.52	252	447317	21.959	ng	100
91) Dibenzo(a,h)anthracene	17.11	278	59555	3.397	ng	95
92) Benzo(g,h,i)perylene	17.56	276	199627	11.431	ng	99

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

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