

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111819\
 Data File : BF117844.D
 Acq On : 18 Nov 2019 19:20
 Operator : JU
 Sample : K5865-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 4-(2-4)

Manual Integrations
 APPROVED

mohammad
 11/19/2019 2:36:07 PM

Quant Time: Nov 19 00:22:49 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	235737	20.00	ng	0.00
21) Naphthalene-d8	8.20	136	897181	20.00	ng	-0.01
39) Acenaphthene-d10	9.97	164	443248	20.00	ng	0.00
64) Phenanthrene-d10	11.46	188	689637	20.00	ng	0.00
76) Chrysene-d12	14.11	240	474406	20.00	ng	0.00
87) Perylene-d12	15.63	264	535880	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	162152	12.18	ng	0.00
7) Phenol-d6	6.53	99	1185926	73.83	ng	-0.01
23) Nitrobenzene-d5	7.48	82	888753	58.89	ng	-0.01
42) 2,4,6-Tribromophenol	10.76	330	19790	4.22	ng	-0.01
45) 2-Fluorobiphenyl	9.29	172	1633450	66.11	ng	0.00
79) Terphenyl-d14	13.05	244	1306566	50.33	ng	0.00
Target Compounds						
10) Phenol	6.55	94	115290	6.907	ng	94
50) Dimethylphthalate	9.67	163	247515	7.963	ng	99
71) Phenanthrene	11.49	178	133901	3.862	ng	95
78) Pyrene	12.91	202	143459	3.742	ng	98
83) Chrysene	14.13	228	62325	2.052	ng	96
88) Benzo(b)fluoranthene	15.17	252	78901m	2.266	ng	

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

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