

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF081018\
 Data File : BF108081.D
 Acq On : 10 Aug 2018 14:00
 Operator : JU/SJ
 Sample : J4276-01RE 5X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-02-080718RE

Manual Integrations
 APPROVED

Sohil
 8/13/2018 3:51:42 PM

Quant Time: Aug 10 15:03:04 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF080818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 08 12:24:24 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.01	152	327064	20.00	ng	0.00
21) Naphthalene-d8	8.30	136	1259225	20.00	ng	0.00
38) Acenaphthene-d10	10.07	164	617411	20.00	ng	0.00
63) Phenanthrene-d10	11.56	188	1037850	20.00	ng	0.00
75) Chrysene-d12	14.22	240	724483	20.00	ng	0.00
86) Perylene-d12	15.78	264	751297	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	168011	9.30	ng	0.00
7) Phenol-d6	6.62	99	238649	9.94	ng	-0.01
23) Nitrobenzene-d5	7.57	82	154729	6.86	ng	0.00
41) 2,4,6-Tribromophenol	10.85	330	43298	7.03	ng	0.00
44) 2-Fluorobiphenyl	9.37	172	319168	8.30	ng	0.00
78) Terphenyl-d14	13.15	244	273764	9.61	ng	0.00
Target Compounds						
57) Fluorene	10.61	166	107712	2.696	ng	98
70) Phenanthrene	11.58	178	897975	18.344	ng	99
71) Anthracene	11.64	178	286735	5.715	ng	100
72) Carbazole	11.78	167	118061	2.649	ng	98
74) Fluoranthene	12.78	202	1099767	20.585	ng	95
77) Pyrene	13.01	202	998782	21.775	ng	99
80) Benzo(a)anthracene	14.21	228	606808	14.594	ng	97
82) Chrysene	14.24	228	489026	12.829	ng	97
85) Indeno(1,2,3-cd)pyrene	17.40	276	234803m	7.109	ng	
87) Benzo(b)fluoranthene	15.31	252	693100m	15.439	ng	
88) Benzo(k)fluoranthene	15.34	252	223121m	5.407	ng	
89) Benzo(a)pyrene	15.71	252	502855	12.509	ng	97
91) Benzo(g,h,i)perylene	17.89	276	199948m	6.105	ng	

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

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