

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082418\
 Data File : BF108485.D
 Acq On : 25 Aug 2018 1:12
 Operator : JU/SJ
 Sample : J4612-01 2X
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RBR200028

Manual Integrations
 APPROVED

Sohil
 8/27/2018 12:56:21 PM

Quant Time: Aug 25 04:07:13 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 22 11:01:45 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.00	152	127448	20.00	ng	0.00
21) Naphthalene-d8	8.28	136	472490	20.00	ng	0.00
38) Acenaphthene-d10	10.05	164	185487	20.00	ng	0.00
63) Phenanthrene-d10	11.55	188	267989	20.00	ng	0.00
75) Chrysene-d12	14.20	240	241112	20.00	ng	0.00
86) Perylene-d12	15.75	264	268343	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	318490	45.24	ng	0.00
7) Phenol-d6	6.62	99	427667	45.72	ng	-0.01
23) Nitrobenzene-d5	7.56	82	272554	32.19	ng	-0.01
41) 2,4,6-Tribromophenol	10.85	330	66871	36.14	ng	0.00
44) 2-Fluorobiphenyl	9.35	172	491209	42.52	ng	-0.01
78) Terphenyl-d14	13.14	244	316209	33.34	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.75	163	52717	4.080	ng	# 80
70) Phenanthrene	11.58	178	178400	14.114	ng	# 95
71) Anthracene	11.63	178	36702	2.833	ng	# 69
74) Fluoranthene	12.78	202	226584	16.425	ng	92
77) Pyrene	13.01	202	191760	12.562	ng	97
80) Benzo(a)anthracene	14.19	228	105925	7.655	ng	95
82) Chrysene	14.22	228	98156m	7.737	ng	
83) Bis(2-ethylhexyl)phthalate	14.15	149	32262	3.930	ng	# 89
85) Indeno(1,2,3-cd)pyrene	17.35	276	44070	4.009	ng	# 85
87) Benzo(b)fluoranthene	15.28	252	127233m	7.935	ng	
88) Benzo(k)fluoranthene	15.31	252	46360m	3.145	ng	
89) Benzo(a)pyrene	15.68	252	89178	6.211	ng	98
91) Benzo(a,h,i)perylene	17.84	276	38563	3.296	ng	# 90

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

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