

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082118\
 Data File : BF108377.D
 Acq On : 22 Aug 2018 9:58
 Operator : JU/SJ
 Sample : J4555-05 2X
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-02-COMP

Manual Integrations
 APPROVED

Sohil
 8/22/2018 4:03:39 PM

Quant Time: Aug 22 11:29:37 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	7.01	152	144864	20.00	ng	0.00
21) Naphthalene-d8	8.29	136	521452	20.00	ng	0.00
38) Acenaphthene-d10	10.05	164	233839	20.00	ng	0.00
63) Phenanthrene-d10	11.55	188	378378	20.00	ng	0.00
75) Chrysene-d12	14.20	240	327317	20.00	ng	0.00
86) Perylene-d12	15.76	264	249902	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	388787	48.59	ng	0.00
7) Phenol-d6	6.62	99	515310	48.46	ng	-0.01
23) Nitrobenzene-d5	7.57	82	323699	34.64	ng	0.00
41) 2,4,6-Tribromophenol	10.84	330	98492	42.23	ng	0.00
44) 2-Fluorobiphenyl	9.36	172	551655	37.87	ng	0.00
78) Terphenyl-d14	13.13	244	431139	33.49	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.75	163	63867	3.921	ng	Qvalue # 99
70) Phenanthrene	11.57	178	103822	5.817	ng	98
74) Fluoranthene	12.77	202	162459	8.341	ng	93
77) Pyrene	13.00	202	137770	6.648	ng	99
80) Benzo(a)anthracene	14.19	228	72165	3.842	ng	97
82) Chrysene	14.22	228	58286	3.384	ng	96
87) Benzo(b)fluoranthene	15.29	252	62862m	4.210	ng	
89) Benzo(a)pyrene	15.69	252	42092	3.148	ng	# 89
91) Benzo(g,h,i)perylene	17.87	276	22864	2.099	ng	# 87

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

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