

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082218\
 Data File : BF108417.D
 Acq On : 23 Aug 2018 8:29
 Operator : JU/SJ
 Sample : J4555-03
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-01-COMP

Manual Integrations
 APPROVED

Sohil
 8/23/2018 3:43:47 PM

Quant Time: Aug 23 09:04:56 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	176445	20.00	ng	0.00
21) Naphthalene-d8	8.28	136	653527	20.00	ng	0.00
38) Acenaphthene-d10	10.04	164	297347	20.00	ng	-0.01
63) Phenanthrene-d10	11.54	188	477945	20.00	ng	0.00
75) Chrysene-d12	14.19	240	334796	20.00	ng	-0.01
86) Perylene-d12	15.75	264	277658	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.63	112	938299	96.28	ng	0.00
7) Phenol-d6	6.62	99	1252427	96.71	ng	0.00
23) Nitrobenzene-d5	7.57	82	810630	69.22	ng	0.00
41) 2,4,6-Tribromophenol	10.84	330	251994	84.96	ng	-0.01
44) 2-Fluorobiphenyl	9.36	172	1306587	70.55	ng	0.00
78) Terphenyl-d14	13.12	244	923755	70.15	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.75	163	178408	8.613	ng	98
70) Phenanthrene	11.56	178	55313	2.454	ng	97
74) Fluoranthene	12.76	202	95962	3.900	ng	92
77) Pyrene	12.99	202	86368	4.075	ng	99
80) Benzo(a)anthracene	14.18	228	48355	2.517	ng	98
82) Chrysene	14.22	228	43493	2.469	ng	98
87) Benzo(b)fluoranthene	15.28	252	49774m	3.000	ng	
88) Benzo(k)fluoranthene	15.28	252	68283	4.477	ng	96
89) Benzo(a)pyrene	15.68	252	34104	2.295	ng	95

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

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