

Data Path : Z:\HPCHEM1\BNA F\DATA\BF022718\
 Data File : BF103266.D
 Acq On : 27 Feb 2018 16:42
 Operator : SJ/JU
 Sample : J1400-11
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-02-022618-C

Manual Integrations
 APPROVED

Sohil
 2/28/2018 11:17:46 AM

Quant Time: Feb 27 17:42:12 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 27 17:05:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	143075	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	576207	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	226338	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	328030	20.00	ng	0.00
75) Chrysene-d12	14.06	240	232645	20.00	ng	0.00
86) Perylene-d12	15.56	264	170672	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	969925	102.53	ng	0.00
7) Phenol-d6	6.51	99	1150395	99.38	ng	0.00
23) Nitrobenzene-d5	7.44	82	717027	71.07	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	171525	89.53	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	956814	69.74	ng	0.00
78) Terphenyl-d14	12.99	244	647659	66.92	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.62	163	90473	4.983	ng	99
70) Phenanthrene	11.43	178	64041	3.547	ng	98
74) Fluoranthene	12.63	202	74317	4.097	ng	97
77) Pyrene	12.86	202	71156	3.923	ng	99
82) Chrysene	14.08	228	29582	2.018	ng	# 92
87) Benzo(b)fluoranthene	15.12	252	26414m	2.354	ng	

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

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