

Data Path : Z:\HPCHEM1\BNA F\DATA\BF032516\
 Data File : BF085776.D
 Acq On : 25 Mar 2016 22:12
 Operator : UM/SJ
 Sample : H1855-12 2X
 Misc :
 ALS Vial : 62 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-7A

Manual Integrations
 APPROVED

Sohil
 3/28/2016 7:59:56 PM

Quant Time: Mar 26 00:13:27 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032416.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 24 14:15:39 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	124204	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	454995	20.00	ng	-0.01
38) Acenaphthene-d10	9.86	164	194244	20.00	ng	0.00
63) Phenanthrene-d10	11.34	188	282952	20.00	ng	-0.01
75) Chrysene-d12	13.98	240	256541	20.00	ng	0.00
86) Perylene-d12	15.40	264	226857	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	26862	3.60	ng	0.01
7) Phenol-d6	6.46	99	36782	3.88	ng	0.00
23) Nitrobenzene-d5	7.38	82	17400	2.15	ng	-0.01
41) 2,4,6-Tribromophenol	10.65	330	4847	2.63	ng	0.00
44) 2-Fluorobiphenyl	9.19	172	30801	2.65	ng	0.00
78) Terphenyl-d14	12.93	244	20037	1.61	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.58	163	120299	8.16	ng	99
70) Phenanthrene	11.36	178	57219	3.94	ng	99
74) Fluoranthene	12.55	202	159308	10.32	ng	97
77) Pyrene	12.78	202	148422	7.04	ng	100
80) Benzo(a)anthracene	13.97	228	97536	6.67	ng	95
82) Chrysene	14.00	228	61840	4.20	ng	98
87) Benzo(b)fluoranthene	15.00	252	111598m	7.81	ng	
88) Benzo(k)fluoranthene	15.02	252	37560m	2.92	ng	
89) Benzo(a)pyrene	15.34	252	76774m	6.05	ng	
91) Benzo(g,h,i)perylene	17.19	276	37654	3.15	ng	98

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

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