

Data Path : Z:\HPCHEM1\BNA F\DATA\BF041718\
 Data File : BF104575.D
 Acq On : 17 Apr 2018 21:29
 Operator : JU/SJ
 Sample : J2387-06
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 20180247-FIELD-DUPLICATE

Manual Integrations
 APPROVED

Sohil
 4/18/2018 3:01:02 PM

Quant Time: Apr 18 02:02:38 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 17 17:08:38 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	140594	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	560301	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	229909	20.00	ng	0.00
63) Phenanthrene-d10	11.33	188	369443	20.00	ng	0.00
75) Chrysene-d12	13.98	240	299325	20.00	ng	0.00
86) Perylene-d12	15.44	264	252711	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	859907	93.52	ng	0.00
7) Phenol-d6	6.45	99	1051011	93.03	ng	0.00
23) Nitrobenzene-d5	7.37	82	634469	73.94	ng	0.00
41) 2,4,6-Tribromophenol	10.63	330	218240	91.51	ng	0.00
44) 2-Fluorobiphenyl	9.16	172	987712	67.87	ng	0.00
78) Terphenyl-d14	12.92	244	788099	60.03	ng	0.00
Target Compounds						
48) Acenaphthylene	9.70	152	54846	2.291	ng	Qvalue 100
49) Dimethylphthalate	9.56	163	99256	5.475	ng	99
70) Phenanthrene	11.36	178	69898	3.397	ng	98
71) Anthracene	11.40	178	52979	2.533	ng	99
74) Fluoranthene	12.54	202	158970	7.395	ng	99
77) Pyrene	12.77	202	208237	9.386	ng	98
80) Benzo(a)anthracene	13.97	228	134507m	7.522	ng	
82) Chrysene	14.00	228	154274	8.399	ng	97
85) Indeno(1,2,3-cd)pyrene	16.88	276	35964	2.305	ng	97
87) Benzo(b)fluoranthene	15.02	252	107801m	6.754	ng	
88) Benzo(k)fluoranthene	15.04	252	32714m	2.171	ng	
89) Benzo(a)pyrene	15.37	252	108011	7.563	ng	95
91) Benzo(a,h,i)perylene	17.32	276	38503m	3.388	ng	

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

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