

Data Path : U:\HPCHEM1\BNA F\DATA\BF020518\
 Data File : BF102756.D
 Acq On : 6 Feb 2018 00:23
 Operator : SJ/JU
 Sample : J1447-05
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 A2 SB45 BOT-S 8.5

Manual Integrations
 APPROVED

Sohil
 2/6/2018 2:25:53 PM

Quant Time: Feb 06 02:48:01 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 05 12:44:16 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.88	152	195783	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	808732	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	350342	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	538895	20.00	ng	0.00
75) Chrysene-d12	14.04	240	324356	20.00	ng	0.00
86) Perylene-d12	15.52	264	311031	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1305262	101.25	ng	0.00
7) Phenol-d6	6.50	99	1522939	98.11	ng	0.00
23) Nitrobenzene-d5	7.44	82	948972	81.40	ng	-0.01
41) 2,4,6-Tribromophenol	10.71	330	279039	98.96	ng	0.00
44) 2-Fluorobiphenyl	9.24	172	1375387	65.14	ng	0.00
78) Terphenyl-d14	12.99	244	877120	64.03	ng	0.00
Target Compounds						
10) Phenol	6.52	94	76661	4.608	ng	98
31) Naphthalene	8.18	128	129525	3.278	ng	99
32) Benzoic acid	7.92	122	4380m	9.072	ng	
49) Dimethylphthalate	9.63	163	223125	8.168	ng	99
70) Phenanthrene	11.43	178	76938	2.563	ng	98
74) Fluoranthene	12.62	202	120142	3.979	ng	99
77) Pyrene	12.85	202	105900	4.164	ng	97
87) Benzo(b)fluoranthene	15.09	252	54392m	2.697	ng	

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

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