

Data Path : U:\HPCHEM1\BNA F\DATA\BF012518\
 Data File : BF102437.D
 Acq On : 25 Jan 2018 21:32
 Operator : SJ/JU
 Sample : J1256-02
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DB-OF-E-W-(SW4-8)

Manual Integrations
 APPROVED

Sohil
 1/26/2018 11:24:38 AM

Quant Time: Jan 26 00:55:08 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 25 13:44:56 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	129028	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	527393	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	231980	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	357709	20.00	ng	0.00
75) Chrysene-d12	13.88	240	206963	20.00	ng	0.00
86) Perylene-d12	15.33	264	194271	20.00	ng	0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	722415	84.89	ng	0.00
7) Phenol-d6	6.36	99	880991	85.87	ng	-0.01
23) Nitrobenzene-d5	7.28	82	532659	58.04	ng	-0.01
41) 2,4,6-Tribromophenol	10.55	330	167915	73.05	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	858284	54.40	ng	0.00
78) Terphenyl-d14	12.83	244	565700	61.62	ng	0.00
Target Compounds						
10) Phenol	6.37	94	43821	3.913	ng	94
49) Dimethylphthalate	9.47	163	147336	7.684	ng	100
70) Phenanthrene	11.26	178	50801	2.437	ng	97
74) Fluoranthene	12.45	202	58353	2.745	ng	95
77) Pyrene	12.68	202	50088m	3.130	ng	
87) Benzo(b)fluoranthene	14.91	252	26195m	2.080	ng	

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

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