

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020818\
 Data File : BF102849.D
 Acq On : 8 Feb 2018 12:09
 Operator : SJ/JU
 Sample : J1469-05
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WABUT-3

Manual Integrations
 APPROVED

Sohil
 2/9/2018 4:33:23 PM

Quant Time: Feb 08 14:15:32 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	209860	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	870944	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	399193	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	659314	20.00	ng	0.00
75) Chrysene-d12	14.04	240	378037	20.00	ng	0.00
86) Perylene-d12	15.53	264	378212	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1374976	99.50	ng	0.01
7) Phenol-d6	6.51	99	1657271	99.60	ng	0.00
23) Nitrobenzene-d5	7.45	82	1101491	87.73	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	318378	99.09	ng	0.00
44) 2-Fluorobiphenyl	9.24	172	1603685	66.66	ng	0.00
78) Terphenyl-d14	13.00	244	1206990	75.60	ng	0.00
Target Compounds						
10) Phenol	6.52	94	40726	2.284	ng	95
32) Benzoic acid	7.88	122	3989m	8.990	ng	
49) Dimethylphthalate	9.63	163	168802	5.423	ng	99
82) Chrysene	14.07	228	57968	2.419	ng	99
87) Benzo(b)fluoranthene	15.09	252	91222m	3.720	ng	

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

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