

Data Path : Z:\HPCHEM1\BNA F\DATA\BF041318\
 Data File : BF104480.D
 Acq On : 13 Apr 2018 16:01
 Operator : JU/SJ
 Sample : J2364-05
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-B

Manual Integrations
 APPROVED

Sohil
 4/16/2018 3:10:12 PM

Quant Time: Apr 16 11:51:19 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 13 17:38:31 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	139928	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	557791	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	246562	20.00	ng	0.00
63) Phenanthrene-d10	11.34	188	379609	20.00	ng	0.00
75) Chrysene-d12	13.99	240	269970	20.00	ng	0.00
86) Perylene-d12	15.45	264	231486	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	869683	95.03	ng	0.00
7) Phenol-d6	6.45	99	1068038	94.99	ng	0.00
23) Nitrobenzene-d5	7.37	82	667807	78.18	ng	0.00
41) 2,4,6-Tribromophenol	10.64	330	235367	92.02	ng	0.00
44) 2-Fluorobiphenyl	9.17	172	1008035	64.59	ng	0.00
78) Terphenyl-d14	12.93	244	764471	64.56	ng	0.00
Target Compounds						
10) Phenol	6.46	94	38962	3.266	ng	95
49) Dimethylphthalate	9.56	163	156383	8.043	ng	100
74) Fluoranthene	12.55	202	63339	2.868	ng	98
77) Pyrene	12.78	202	60153	3.006	ng	97
82) Chrysene	14.01	228	33965	2.050	ng	98
87) Benzo(b)fluoranthene	15.02	252	38339m	2.622	ng	

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

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