

Data Path : Z:\HPCHEM1\BNA F\DATA\BF030218\
 Data File : BF103364.D
 Acq On : 2 Mar 2018 14:10
 Operator : SJ/JU
 Sample : J1402-13
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PL-01-022818-D

Manual Integrations
 APPROVED

Sohil
 3/5/2018 3:13:08 PM

Quant Time: Mar 05 10:23:18 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 02 13:28:09 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	167564	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	657233	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	269523	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	401178	20.00	ng	0.00
75) Chrysene-d12	14.06	240	263096	20.00	ng	0.00
86) Perylene-d12	15.57	264	236874	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1093027	98.66	ng	0.00
7) Phenol-d6	6.52	99	1318453	97.25	ng	0.00
23) Nitrobenzene-d5	7.44	82	800138	69.53	ng	-0.01
41) 2,4,6-Tribromophenol	10.71	330	187945	82.38	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	958459	56.39	ng	0.00
78) Terphenyl-d14	12.99	244	575331	52.57	ng	0.00
Target Compounds						
10) Phenol	6.52	94	55487	3.526	ng	Qvalue # 72
49) Dimethylphthalate	9.63	163	92151	4.262	ng	99
70) Phenanthrene	11.43	178	64670	2.929	ng	97
74) Fluoranthene	12.63	202	137416	6.194	ng	99
77) Pyrene	12.86	202	119561	5.829	ng	98
80) Benzo(a)anthracene	14.05	228	60214	3.527	ng	97
82) Chrysene	14.09	228	70622	4.261	ng	98
85) Indeno(1,2,3-cd)pyrene	17.10	276	30940	2.358	ng	98
87) Benzo(b)fluoranthene	15.13	252	92651m	5.949	ng	
89) Benzo(a)pyrene	15.50	252	56210	4.042	ng	# 92
91) Benzo(g,h,i)perylene	17.57	276	33377	3.178	ng	# 83

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

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