

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030518\
 Data File : BF103405.D
 Acq On : 5 Mar 2018 13:17
 Operator : SJ/JU
 Sample : J1751-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-02-030118-A

Manual Integrations
 APPROVED

Sohil
 3/6/2018 4:22:45 PM

Quant Time: Mar 05 14:01:58 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 05 13:40:47 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	138195	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	564286	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	236248	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	337390	20.00	ng	0.00
75) Chrysene-d12	14.06	240	244036	20.00	ng	0.00
86) Perylene-d12	15.58	264	208820	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	840231	91.96	ng	0.00
7) Phenol-d6	6.52	99	1028229	91.96	ng	0.00
23) Nitrobenzene-d5	7.44	82	625180	63.27	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	143390	71.71	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	853760	57.50	ng	0.00
78) Terphenyl-d14	12.99	244	543287	53.52	ng	0.00
Target Compounds						
10) Phenol	6.53	94	41977	3.234	ng	Qvalue # 59
49) Dimethylphthalate	9.62	163	271932	14.348	ng	99
70) Phenanthrene	11.43	178	67720	3.647	ng	99
74) Fluoranthene	12.63	202	118297	6.340	ng	97
77) Pyrene	12.86	202	111012	5.835	ng	98
80) Benzo(a)anthracene	14.05	228	52236	3.299	ng	96
82) Chrysene	14.09	228	47298m	3.076	ng	
87) Benzo(b)fluoranthene	15.13	252	53400m	3.889	ng	
89) Benzo(a)pyrene	15.52	252	36368	2.966	ng	# 79

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

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