

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF031618\
 Data File : BF103702.D
 Acq On : 16 Mar 2018 17:59
 Operator : JU/SJ
 Sample : J1752-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NB-08-031418

Manual Integrations
 APPROVED

Sohil
 3/19/2018 2:00:57 PM

Quant Time: Mar 19 10:37:34 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 15 18:31:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	180931	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	740057	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	337498	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	570316	20.00	ng	0.00
75) Chrysene-d12	14.14	240	374950	20.00	ng	0.00
86) Perylene-d12	15.68	264	354855	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	1197305	100.65	ng	0.00
7) Phenol-d6	6.59	99	1468550	100.91	ng	0.00
23) Nitrobenzene-d5	7.53	82	891794	84.03	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	305795	105.38	ng	0.00
44) 2-Fluorobiphenyl	9.32	172	1362552	64.40	ng	0.00
78) Terphenyl-d14	13.09	244	1112730	68.88	ng	0.00
Target Compounds						
10) Phenol	6.59	94	43329	2.802	ng	Qvalue # 54
32) Benzoic acid	7.91	122	3411m	10.270	ng	
49) Dimethylphthalate	9.71	163	131427	5.104	ng	99
74) Fluoranthene	12.71	202	146954	4.572	ng	97
77) Pyrene	12.94	202	124146	4.508	ng	98
80) Benzo(a)anthracene	14.13	228	52445	2.210	ng	95
82) Chrysene	14.17	228	56693m	2.506	ng	
87) Benzo(b)fluoranthene	15.22	252	68410	3.051	ng	# 85
89) Benzo(a)pyrene	15.61	252	45695	2.247	ng	# 89

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

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