

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF031618\
 Data File : BF103703.D
 Acq On : 16 Mar 2018 18:26
 Operator : JU/SJ
 Sample : J1757-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-01-031418-A

Manual Integrations
 APPROVED

Sohil
 3/19/2018 2:01:00 PM

Quant Time: Mar 17 00:44:55 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	6.97	152	173476	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	689354	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	306598	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	491553	20.00	ng	0.00
75) Chrysene-d12	14.15	240	325146	20.00	ng	0.00
86) Perylene-d12	15.69	264	304386	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	1106528	97.02	ng	0.00
7) Phenol-d6	6.59	99	1371742	98.31	ng	0.00
23) Nitrobenzene-d5	7.53	82	819515	82.90	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	277261	105.18	ng	0.00
44) 2-Fluorobiphenyl	9.32	172	1161034	60.41	ng	0.00
78) Terphenyl-d14	13.09	244	898408	64.14	ng	0.00
Target Compounds						
10) Phenol	6.60	94	43346	2.923	ng	95
49) Dimethylphthalate	9.71	163	103293	4.416	ng	# 99
70) Phenanthrene	11.52	178	65616	2.440	ng	98
74) Fluoranthene	12.72	202	74218	2.679	ng	96
77) Pyrene	12.95	202	89831	3.762	ng	99
82) Chrysene	14.17	228	40889	2.084	ng	# 94
87) Benzo(b)fluoranthene	15.22	252	48197m	2.506	ng	
89) Benzo(a)pyrene	15.61	252	35748	2.050	ng	# 65

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

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