

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF021618\
 Data File : BF103062.D
 Acq On : 16 Feb 2018 19:39
 Operator : SJ/JU
 Sample : J1584-03
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-GF-02-012-B

Manual Integrations
 APPROVED

Sohil
 2/19/2018 11:09:53 AM

Quant Time: Feb 16 23:44:09 2018
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 16 14:34:49 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	171503	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	662271	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	250753	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	378352	20.00	ng	0.00
75) Chrysene-d12	14.04	240	286908	20.00	ng	-0.01
86) Perylene-d12	15.54	264	232634	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1111757	98.45	ng	0.01
7) Phenol-d6	6.51	99	1314444	96.67	ng	0.00
23) Nitrobenzene-d5	7.44	82	801348	83.94	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	194601	96.42	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1118816	74.04	ng	0.00
78) Terphenyl-d14	12.99	244	854089	70.49	ng	0.00
Target Compounds						
10) Phenol	6.52	94	34296	2.354	ng	Qvalue 89
49) Dimethylphthalate	9.62	163	93977	4.807	ng	100
70) Phenanthrene	11.43	178	50358	2.390	ng	97
74) Fluoranthene	12.62	202	92528	4.364	ng	96
77) Pyrene	12.85	202	89278	3.968	ng	98
80) Benzo(a)anthracene	14.04	228	41711	2.312	ng	93
82) Chrysene	14.07	228	42414	2.332	ng	96
87) Benzo(b)fluoranthene	15.10	252	44426m	2.946	ng	
89) Benzo(a)pyrene	15.47	252	28919	2.130	ng	# 91

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

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