

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112818\
 Data File : BF111055.D
 Acq On : 28 Nov 2018 19:43
 Operator : JU/SJ
 Sample : J6080-03
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RCA-06

Manual Integrations
 APPROVED

mohammad
 11/29/2018 1:58:30 PM

Quant Time: Nov 29 03:05:24 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF112618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 27 15:32:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	59021	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	239467	20.00	ng	0.00
39) Acenaphthene-d10	9.97	164	98795	20.00	ng	0.00
64) Phenanthrene-d10	11.47	188	149578	20.00	ng	0.00
76) Chrysene-d12	14.13	240	120679	20.00	ng	0.00
87) Perylene-d12	15.66	264	98122	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	318132	89.39	ng	0.00
7) Phenol-d6	6.56	99	404215	87.15	ng	0.00
23) Nitrobenzene-d5	7.50	82	261465	62.49	ng	-0.01
42) 2,4,6-Tribromophenol	10.77	330	76543	78.01	ng	-0.01
45) 2-Fluorobiphenyl	9.29	172	409356	60.03	ng	0.00
79) Terphenyl-d14	13.05	244	279683	45.11	ng	0.00
Target Compounds						
10) Phenol	6.57	94	23886	4.573	ng	83
50) Dimethylphthalate	9.69	163	27464	3.790	ng	# 99
71) Phenanthrene	11.50	178	51883	6.549	ng	99
75) Fluoranthene	12.69	202	57620	7.077	ng	99
78) Pyrene	12.93	202	50652	5.747	ng	97
81) Benzo(a)anthracene	14.12	228	23221	3.259	ng	# 96
83) Chrysene	14.16	228	21836	3.113	ng	95
88) Benzo(b)fluoranthene	15.22	252	20481m	3.553	ng	
90) Benzo(a)pyrene	15.60	252	15852	2.954	ng	# 94

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

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