

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062818\
 Data File : BF106941.D
 Acq On : 29 Jun 2018 00:30
 Operator : JU/SJ
 Sample : J3048-15
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-2-D

Manual Integrations
 APPROVED

Sohil
 6/29/2018 2:30:57 PM

Quant Time: Jun 29 03:28:05 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 25 18:50:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	61763	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	211420	20.00	ng	-0.01
38) Acenaphthene-d10	10.00	164	86259	20.00	ng	-0.01
63) Phenanthrene-d10	11.48	188	148003	20.00	ng	-0.02
75) Chrysene-d12	14.14	240	197668	20.00	ng	-0.03
86) Perylene-d12	15.67	264	182308	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	104859	28.75	ng	0.00
7) Phenol-d6	6.60	99	355356	78.02	ng	-0.02
23) Nitrobenzene-d5	7.52	82	320392	84.22	ng	-0.02
41) 2,4,6-Tribromophenol	10.79	330	13984	13.99	ng	-0.01
44) 2-Fluorobiphenyl	9.31	172	402320	77.67	ng	-0.02
78) Terphenyl-d14	13.07	244	450329	55.19	ng	-0.01
Target Compounds						
10) Phenol	6.61	94	18958	3.647	ng	99
20) 3+4-Methylphenols	7.36	107	40996	8.549	ng	88
32) Benzoic acid	8.02	122	455	5.925	ng	# 1
49) Dimethylphthalate	9.70	163	130943	20.240	ng	99
70) Phenanthrene	11.51	178	33503	4.693	ng	96
74) Fluoranthene	12.70	202	71452	9.081	ng	96
77) Pyrene	12.94	202	68365	5.245	ng	97
80) Benzo(a)anthracene	14.12	228	49850	4.138	ng	97
82) Chrysene	14.16	228	44451	4.011	ng	97
87) Benzo(b)fluoranthene	15.21	252	54177m	4.784	ng	
89) Benzo(a)pyrene	15.60	252	38084	3.783	ng	97

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

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