

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF061918\
 Data File : BF106594.D
 Acq On : 19 Jun 2018 20:43
 Operator : JU/SJ
 Sample : J3528-03
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TPTW-01

Quant Time: Jun 27 09:34:42 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 19 14:32:09 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	112489	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	430084	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	191471	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	317777	20.00	ng	0.00
75) Chrysene-d12	14.15	240	256861	20.00	ng	0.00
86) Perylene-d12	15.69	264	294851	20.00	ng	0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.61	112	385425	58.02	ng	0.00
7) Phenol-d6	6.60	99	305520	36.83	ng	0.00
23) Nitrobenzene-d5	7.52	82	646026	83.48	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	261708	117.93	ng	0.00
44) 2-Fluorobiphenyl	9.32	172	1067815	96.37	ng	0.00
78) Terphenyl-d14	13.08	244	856366	80.76	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
27) 2,4-Dimethylphenol	7.90	122	22631	3.925	ng	92
37) 2-Methylnaphthalene	8.95	142	38194	2.866	ng	98
49) Dimethylphthalate	9.71	163	137463	9.572	ng	99
51) Acenaphthene	10.04	154	51763	4.335	ng	99
57) Fluorene	10.55	166	46404	3.576	ng	97
70) Phenanthrene	11.52	178	90443	5.901	ng	97
72) Carbazole	11.73	167	59977	4.095	ng	97

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

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