

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121818\
 Data File : BM018273.D
 Acq On : 18 Dec 2018 16:23
 Operator : JU/SJ
 Sample : J6403-61
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TP-8-B

Quant Time: Dec 18 17:00:47 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM121518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Dec 15 21:39:37 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.49	152	134064	20.00	ng	-0.02
21) Naphthalene-d8	10.26	136	525411	20.00	ng	-0.01
39) Acenaphthene-d10	14.15	164	282270	20.00	ng	-0.02
64) Phenanthrene-d10	16.91	188	554705	20.00	ng	-0.01
76) Chrysene-d12	21.13	240	459411	20.00	ng	-0.01
87) Perylene-d12	23.30	264	478001	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.12	112	832364	103.72	ng	0.00
7) Phenol-d6	6.69	99	1136698	101.90	ng	-0.01
23) Nitrobenzene-d5	8.65	82	720150	70.31	ng	-0.02
42) 2,4,6-Tribromophenol	15.66	330	356464	86.04	ng	-0.01
45) 2-Fluorobiphenyl	12.75	172	1247881	57.26	ng	-0.02
79) Terphenyl-d14	19.56	244	1419621	69.89	ng	-0.01
Target Compounds						
10) Phenol	6.72	94	57346	4.854	ng	88
50) Dimethylphthalate	13.62	163	177934	7.582	ng	100
75) Fluoranthene	18.99	202	108204	3.089	ng	99
78) Pyrene	19.35	202	96357	3.520	ng	96
88) Benzo(b)fluoranthene	22.66	252	76256	2.840	ng	# 94
90) Benzo(a)pyrene	23.20	252	53387	2.091	ng	99

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

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