

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122118\
 Data File : BM018385.D
 Acq On : 22 Dec 2018 16:11
 Operator : JU/SJ
 Sample : J6494-01
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 OWS

Manual Integrations
 APPROVED

Sohil
 12/26/2018 3:20:33 PM

Quant Time: Dec 24 06:18:20 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	7.47	152	96795	20.00	ng	-0.03
21) Naphthalene-d8	10.24	136	384627	20.00	ng	-0.03
39) Acenaphthene-d10	14.13	164	207259	20.00	ng	-0.03
64) Phenanthrene-d10	16.90	188	391001	20.00	ng	-0.02
76) Chrysene-d12	21.13	240	335428	20.00	ng	-0.02
87) Perylene-d12	23.29	264	381842	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.12	112	919791	158.74	ng	0.00
7) Phenol-d6	6.70	99	1122535	139.38	ng	0.00
23) Nitrobenzene-d5	8.64	82	902500	120.36	ng	-0.02
42) 2,4,6-Tribromophenol	15.65	330	430480	141.51	ng	-0.02
45) 2-Fluorobiphenyl	12.74	172	1748886	109.29	ng	-0.03
79) Terphenyl-d14	19.56	244	1684125	113.55	ng	-0.01
Target Compounds						
10) Phenol	6.72	94	897432	105.217	ng	91
15) Benzyl Alcohol	7.73	79	149038	22.711	ng	92
20) 3+4-Methylphenols	8.26	107	267378	33.505	ng	91
32) Benzoic acid	9.72	122	164645m	32.830	ng	
50) Dimethylphthalate	13.61	163	74500	4.324	ng	98

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

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