

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082718\
 Data File : BM016613.D
 Acq On : 27 Aug 2018 19:04
 Operator : SJ/JU
 Sample : J4652-04
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SONO-W

Manual Integrations
 APPROVED

Sohil
 8/28/2018 11:16:57 AM

Quant Time: Aug 28 10:43:12 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 27 13:01:22 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.00	152	104455	20.00	ng	0.00
21) Naphthalene-d8	10.81	136	385779	20.00	ng	0.00
38) Acenaphthene-d10	14.65	164	320539	20.00	ng	0.00
63) Phenanthrene-d10	17.40	188	1061163	20.00	ng	0.02
75) Chrysene-d12	21.53	240	1334688	20.00	ng	0.00
86) Perylene-d12	23.81	264	1371006	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.55	112	292198	54.82	ng	0.00
7) Phenol-d6	7.19	99	231734	32.85	ng	0.00
23) Nitrobenzene-d5	9.19	82	626384	74.14	ng	0.00
41) 2,4,6-Tribromophenol	16.14	330	1222745	136.46	ng	0.00
44) 2-Fluorobiphenyl	13.26	172	2398452	71.92	ng	0.00
78) Terphenyl-d14	19.98	244	6530398	103.56	ng	0.00
Target Compounds						
10) Phenol	7.21	94	21551	3.089	ng	# 64
42) 2,4,6-Trichlorophenol	13.10	196	29590	2.733	ng	# 93
49) Dimethylphthalate	14.11	163	78053	2.535	ng	# 93
58) 2,3,4,6-Tetrachlorophenol	15.29	232	4751973	467.549	ng	# 100
69) Pentachlorophenol	17.05	266	42147884m	4189.291	ng	

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

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