

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122419\
 Data File : BP001409.D
 Acq On : 25 Dec 2019 06:19
 Operator : JU
 Sample : K6401-04DL 2X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0B38DL

Quant Time: Dec 25 08:21:37 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP121919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 24 16:04:49 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.28	152	39011	20.00	ng	0.00
21) Naphthalene-d8	10.32	136	146251	20.00	ng	0.00
39) Acenaphthene-d10	13.18	164	80546	20.00	ng	0.00
64) Phenanthrene-d10	15.58	188	147583	20.00	ng	0.00
76) Chrysene-d12	19.23	240	116658	20.00	ng	0.00
87) Perylene-d12	21.00	264	123347	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	6.19	112	98575	40.84	ng	0.00
7) Phenol-d6	7.74	99	135330	42.97	ng	-0.01
23) Nitrobenzene-d5	9.16	82	92764	24.36	ng	-0.01
42) 2,4,6-Tribromophenol	14.47	330	35564	35.32	ng	0.00
45) 2-Fluorobiphenyl	12.10	172	174586	27.41	ng	0.00
79) Terphenyl-d14	17.87	244	190474	27.97	ng	0.00
Target Compounds						
17) 2-Methylphenol	8.71	107	29924	13.955	ng	98
20) 3+4-Methylphenols	8.97	107	150408	52.857	ng	94
27) 2,4-Dimethylphenol	9.81	122	50647	25.763	ng	96
50) Dimethylphthalate	12.77	163	25787	3.902	ng	97

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

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