

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122319\
 Data File : BP001368.D
 Acq On : 23 Dec 2019 15:38
 Operator : JU
 Sample : K6372-05
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MW-6

Manual Integrations
 APPROVED

mohammad
 12/24/2019 3:05:22 PM

Quant Time: Jan 01 06:25:56 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP121919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 20 03:20:39 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.30	152	45501	20.00	ng	0.00
21) Naphthalene-d8	10.34	136	162329	20.00	ng	0.00
39) Acenaphthene-d10	13.20	164	96306	20.00	ng	0.00
64) Phenanthrene-d10	15.60	188	190243	20.00	ng	0.00
76) Chrysene-d12	19.25	240	150423	20.00	ng	0.00
87) Perylene-d12	21.02	264	158033	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	6.22	112	170763	60.65	ng	0.02
7) Phenol-d6	7.76	99	110255	30.01	ng	0.00
23) Nitrobenzene-d5	9.19	82	358069	84.71	ng	0.00
42) 2,4,6-Tribromophenol	14.50	330	164678	136.77	ng	0.00
45) 2-Fluorobiphenyl	12.12	172	635604	83.45	ng	0.00
79) Terphenyl-d14	17.89	244	823378	93.75	ng	0.00
Target Compounds						
10) Phenol	7.78	94	21214	5.054	ng	Qvalue # 80
17) 2-Methylphenol	8.73	107	13559	5.421	ng	# 66
27) 2,4-Dimethylphenol	9.85	122	51244m	23.485	ng	
31) Naphthalene	10.38	128	518396	58.503	ng	99
37) 2-Methylnaphthalene	11.50	142	132905	21.726	ng	97
38) 1-Methylnaphthalene	11.66	142	97219m	16.904	ng	

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

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