

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP060220\
 Data File : BP002357.D
 Acq On : 02 Jun 2020 16:01
 Operator : CG/JU
 Sample : L2822-03DL 10X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AP8DL

Manual Integrations
 APPROVED

mohammad
 6/3/2020 4:36:55 PM

Quant Time: Jun 03 02:29:54 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP052720.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 02 12:11:53 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	217114	20.00	ng	0.00
21) Naphthalene-d8	10.38	136	847849	20.00	ng	0.00
39) Acenaphthene-d10	14.25	164	493954	20.00	ng	-0.01
64) Phenanthrene-d10	17.01	188	1040272	20.00	ng	-0.01
76) Chrysene-d12	21.12	240	1063102	20.00	ng	-0.01
86) Perylene-d12	23.33	264	1147716	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.23	112	165410	14.31	ng	0.00
7) Phenol-d6	6.80	99	205107	13.04	ng	0.00
23) Nitrobenzene-d5	8.75	82	161170	11.43	ng	0.00
42) 2,4,6-Tribromophenol	15.75	330	89544	19.13	ng	-0.01
45) 2-Fluorobiphenyl	12.87	172	363862	13.00	ng	-0.01
79) Terphenyl-d14	19.60	244	607845	13.85	ng	0.00
Target Compounds						
10) Phenol	6.82	94	404677	23.387	ng	99
17) 2-Methylphenol	8.06	107	519663	42.047	ng	99
20) 3+4-Methylphenols	8.39	107	1109225	66.046	ng	99
27) 2,4-Dimethylphenol	9.58	122	790981m	73.347	ng	
32) Benzoic acid	9.67	122	20247	6.985	ng	91
53) 3-Nitroaniline	14.16	138	37120m	4.471	ng	

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

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