

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122419\
 Data File : BP001400.D
 Acq On : 24 Dec 2019 22:04
 Operator : JU
 Sample : K6401-06
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0B40

Manual Integrations
 APPROVED

mohammad
 12/26/2019 4:01:40 PM

Quant Time: Dec 30 07:25:59 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	8.32	152	52210	20.00	ng	0.03
21) Naphthalene-d8	10.33	136	186757	20.00	ng	0.01
39) Acenaphthene-d10	13.19	164	84685	20.00	ng	0.01
64) Phenanthrene-d10	15.60	188	163342	20.00	ng	0.02
76) Chrysene-d12	19.23	240	123652	20.00	ng	0.00
87) Perylene-d12	21.02	264	127675	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	6.21	112	272300	84.29	ng	0.01
7) Phenol-d6	7.82	99	351435m	83.37	ng	0.07
23) Nitrobenzene-d5	9.18	82	242289	49.82	ng	0.00
42) 2,4,6-Tribromophenol	14.53	330	92343	87.22	ng	0.05
45) 2-Fluorobiphenyl	12.11	172	417133	62.28	ng	0.01
79) Terphenyl-d14	17.87	244	433447	60.04	ng	0.00
Target Compounds						
9) Benzaldehyde	7.58	77	9657	3.266	ng	Qvalue # 44
12) 1,3-Dichlorobenzene	8.23	146	2772468	685.363	ng	97
13) 1,4-Dichlorobenzene	8.36	146	4183955	1015.182	ng	100
14) 1,2-Dichlorobenzene	8.63	146	7695476	1959.969	ng	99
17) 2-Methylphenol	8.76	107	6998m	2.438	ng	
20) 3+4-Methylphenols	9.00	107	86548	22.726	ng	94
27) 2,4-Dimethylphenol	9.86	122	64317m	25.621	ng	
30) 1,2,4-Trichlorobenzene	10.25	180	370029	96.779	ng	98
50) Dimethylphthalate	12.79	163	81140	11.679	ng	97

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

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