

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP012720\
 Data File : BP001698.D
 Acq On : 27 Jan 2020 19:12
 Operator : CG/JU
 Sample : L1242-14
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 20200067-AMENDED-COMP-5

Manual Integrations
 APPROVED

mohammad
 1/28/2020 1:14:32 PM

Quant Time: Jan 28 04:27:27 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP010820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 24 22:23:46 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	18271	20.00	ng	-0.01
21) Naphthalene-d8	10.33	136	83121	20.00	ng	-0.01
39) Acenaphthene-d10	14.21	164	63939	20.00	ng	-0.01
64) Phenanthrene-d10	16.97	188	148464	20.00	ng	-0.01
76) Chrysene-d12	21.08	240	148376	20.00	ng	-0.01
87) Perylene-d12	23.29	264	180164	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.18	112	6219	6.67	ng	0.00
7) Phenol-d6	6.75	99	33797	22.67	ng	-0.01
23) Nitrobenzene-d5	8.70	82	42195	22.18	ng	-0.01
42) 2,4,6-Tribromophenol	15.72	330	7490	7.70	ng	-0.01
45) 2-Fluorobiphenyl	12.82	172	136912	31.48	ng	-0.02
79) Terphenyl-d14	19.56	244	302823	37.33	ng	-0.02
Target Compounds						
10) Phenol	6.78	94	3704	2.402	ng	95
50) Dimethylphthalate	13.67	163	26080	4.952	ng	98
75) Fluoranthene	19.00	202	35066	3.301	ng	98
78) Pyrene	19.35	202	43501	3.991	ng	98
81) Benzo(a)anthracene	21.07	228	24722	2.392	ng	# 87
83) Chrysene	21.12	228	26687	2.650	ng	93
88) Benzo(b)fluoranthene	22.63	252	27682m	2.438	ng	
90) Benzo(a)pyrene	23.18	252	25834	2.397	ng	# 98

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

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