

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP012720\
 Data File : BP001696.D
 Acq On : 27 Jan 2020 18:04
 Operator : CG/JU
 Sample : L1242-12
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 20200065-AMENDED-COMP-3

Manual Integrations
 APPROVED

mohammad
 1/28/2020 1:14:30 PM

Quant Time: Jan 28 04:21:29 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	20676	20.00	ng	-0.01
21) Naphthalene-d8	10.33	136	93318	20.00	ng	-0.01
39) Acenaphthene-d10	14.21	164	70251	20.00	ng	-0.01
64) Phenanthrene-d10	16.97	188	166961	20.00	ng	-0.01
76) Chrysene-d12	21.08	240	164870	20.00	ng	-0.01
87) Perylene-d12	23.28	264	185940	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.18	112	47386	44.89	ng	0.00
7) Phenol-d6	6.75	99	141463	83.86	ng	-0.01
23) Nitrobenzene-d5	8.70	82	171573	80.34	ng	-0.01
42) 2,4,6-Tribromophenol	15.72	330	22851	21.37	ng	-0.01
45) 2-Fluorobiphenyl	12.83	172	359766	75.29	ng	-0.01
79) Terphenyl-d14	19.57	244	667971	74.10	ng	-0.01
Target Compounds						
10) Phenol	6.78	94	9252	5.302	ng	Qvalue 84
20) 3+4-Methylphenols	8.33	107	10954	6.464	ng	91
50) Dimethylphthalate	13.67	163	48197	8.330	ng	99
71) Phenanthrene	17.02	178	31046	3.419	ng	99
72) Anthracene	17.10	178	18180	2.057	ng	95
75) Fluoranthene	19.00	202	81270	6.802	ng	98
78) Pyrene	19.35	202	70311	5.806	ng	99
81) Benzo(a)anthracene	21.07	228	37528	3.268	ng	91
83) Chrysene	21.12	228	47863	4.277	ng	98
88) Benzo(b)fluoranthene	22.62	252	40771	3.480	ng	# 98
90) Benzo(a)pyrene	23.18	252	30484	2.740	ng	# 96

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

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