

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP012720\
 Data File : BP001695.D
 Acq On : 27 Jan 2020 17:31
 Operator : CG/JU
 Sample : L1242-11
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 20200064-AMENDED-COMP-2

Manual Integrations
 APPROVED

mohammad
 1/28/2020 1:14:27 PM

Quant Time: Jan 28 04:16:48 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	16597	20.00	ng	-0.01
21) Naphthalene-d8	10.33	136	74053	20.00	ng	-0.01
39) Acenaphthene-d10	14.21	164	59677	20.00	ng	-0.01
64) Phenanthrene-d10	16.97	188	142917	20.00	ng	-0.02
76) Chrysene-d12	21.08	240	145154	20.00	ng	-0.01
87) Perylene-d12	23.28	264	173264	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.18	112	52107	61.50	ng	0.00
7) Phenol-d6	6.75	99	110411	81.54	ng	-0.01
23) Nitrobenzene-d5	8.70	82	111314	65.68	ng	-0.01
42) 2,4,6-Tribromophenol	15.72	330	35258	38.81	ng	-0.01
45) 2-Fluorobiphenyl	12.82	172	243606	60.02	ng	-0.02
79) Terphenyl-d14	19.56	244	470877	59.33	ng	-0.02
Target Compounds						
10) Phenol	6.78	94	7040	5.026	ng	86
50) Dimethylphthalate	13.67	163	30426	6.190	ng	99
75) Fluoranthene	19.00	202	31285	3.059	ng	99
78) Pyrene	19.35	202	36998	3.470	ng	97
88) Benzo(b)fluoranthene	22.62	252	22391	2.051	ng	# 96

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

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