

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP090120\
 Data File : BP003169.D
 Acq On : 01 Sep 2020 16:25
 Operator : CG/JU
 Sample : L3848-13
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 LINDEN-BH-04-B

Manual Integrations
 APPROVED

mohammad
 9/2/2020 1:58:46 PM

Quant Time: Sep 08 08:35:55 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP082420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 01 14:37:05 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	124095	20.00	ng	0.00
21) Naphthalene-d8	10.43	136	503281	20.00	ng	0.00
39) Acenaphthene-d10	14.29	164	313234	20.00	ng	0.00
64) Phenanthrene-d10	17.05	188	712838	20.00	ng	0.00
76) Chrysene-d12	21.14	240	958782	20.00	ng	0.00
86) Perylene-d12	23.37	264	1286177	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	487192	69.41	ng	0.00
7) Phenol-d6	6.83	99	630845	71.70	ng	0.00
23) Nitrobenzene-d5	8.80	82	359938	51.56	ng	0.00
42) 2,4,6-Tribromophenol	15.79	330	267303	63.10	ng	0.00
45) 2-Fluorobiphenyl	12.92	172	1102455	51.54	ng	0.00
79) Terphenyl-d14	19.62	244	1913058	43.91	ng	0.00
Target Compounds						
10) Phenol	6.86	94	50148	6.150	ng	95
20) 3+4-Methylphenols	8.43	107	24268	3.101	ng	87
31) Naphthalene	10.48	128	90012	3.481	ng	99
50) Dimethylphthalate	13.75	163	188557	7.894	ng	100
71) Phenanthrene	17.09	178	206202	5.338	ng	99
75) Fluoranthene	19.06	202	247995	5.555	ng	98
78) Pyrene	19.42	202	234670	3.771	ng	98
81) Benzo(a)anthracene	21.12	228	157954	2.630	ng	99
83) Chrysene	21.17	228	146514	2.548	ng	96
88) Benzo(b)fluoranthene	22.70	252	231885m	2.901	ng	
90) Benzo(a)pyrene	23.27	252	183551	2.475	ng	96
92) Benzo(a,h,i)perylene	26.32	276	175006	2.213	ng	97

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

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