

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP090220\
 Data File : BP003183.D
 Acq On : 02 Sep 2020 13:04
 Operator : CG/JU
 Sample : L3848-09RE
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 LINDEN-BH-01-ARE

Manual Integrations
 APPROVED

mohammad
 9/3/2020 1:48:18 PM

Quant Time: Sep 02 15:15:19 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	111011	20.00	ng	0.00
21) Naphthalene-d8	10.43	136	470853	20.00	ng	0.00
39) Acenaphthene-d10	14.29	164	310539	20.00	ng	0.00
64) Phenanthrene-d10	17.04	188	743956	20.00	ng	0.00
76) Chrysene-d12	21.14	240	1118967	20.00	ng	0.00
86) Perylene-d12	23.37	264	1398177	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	1361	0.22	ng	0.00
7) Phenol-d6	6.83	99	11906	1.51	ng	0.00
23) Nitrobenzene-d5	8.79	82	275207	42.14	ng	0.00
42) 2,4,6-Tribromophenol	15.79	330	1032	0.25	ng	0.00
45) 2-Fluorobiphenyl	12.91	172	842712	39.74	ng	0.00
79) Terphenyl-d14	19.62	244	1684038	33.12	ng	0.00
Target Compounds						
25) Isophorone	9.36	82	28157	2.443	ng	Qvalue 96
75) Fluoranthene	19.07	202	225395	4.838	ng	100
78) Pyrene	19.42	202	247451	3.407	ng	99
81) Benzo(a)anthracene	21.13	228	185747	2.650	ng	95
83) Chrysene	21.17	228	191755m	2.857	ng	
88) Benzo(b)fluoranthene	22.70	252	302656m	3.483	ng	
90) Benzo(a)pyrene	23.27	252	216912	2.691	ng	# 93

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

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