

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP050520\
 Data File : BP002245.D
 Acq On : 05 May 2020 19:52
 Operator : CG/JU
 Sample : L2483-04
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 6FT-TP-2

Manual Integrations
 APPROVED

mohammad
 5/6/2020 6:44:35 PM

Quant Time: May 06 01:33:05 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP050420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 05 13:34:25 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	271738	20.00	ng	0.00
21) Naphthalene-d8	10.39	136	1230641	20.00	ng	0.00
39) Acenaphthene-d10	14.26	164	781549	20.00	ng	0.00
64) Phenanthrene-d10	17.01	188	1705220	20.00	ng	0.00
76) Chrysene-d12	21.11	240	1749466	20.00	ng	0.00
87) Perylene-d12	23.31	264	1808985	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.23	112	1474510	89.52	ng	0.00
7) Phenol-d6	6.79	99	2243771	92.73	ng	0.00
23) Nitrobenzene-d5	8.76	82	1499369	62.92	ng	0.00
42) 2,4,6-Tribromophenol	15.75	330	816433	98.70	ng	0.00
45) 2-Fluorobiphenyl	12.87	172	3363555	60.27	ng	0.00
79) Terphenyl-d14	19.60	244	4825811	67.60	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.72	163	649897	10.550	ng	99
71) Phenanthrene	17.05	178	195415	2.038	ng	99
75) Fluoranthene	19.03	202	442378	4.021	ng	96
78) Pyrene	19.38	202	390901	3.289	ng	98
83) Chrysene	21.14	228	227472m	2.041	ng	
88) Benzo(b)fluoranthene	22.65	252	309407	2.844	ng	97
90) Benzo(a)pyrene	23.21	252	211415	2.019	ng	96

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

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