

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081920\
 Data File : BP003045.D
 Acq On : 19 Aug 2020 22:20
 Operator : CG/JU
 Sample : L3712-05
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 VB-25203

Manual Integrations
 APPROVED

mohammad
 8/20/2020 12:24:42 PM

Quant Time: Aug 20 02:32:52 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP073020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 17 14:07:35 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	342468	20.00	ng	0.00
21) Naphthalene-d8	10.46	136	1397156	20.00	ng	-0.01
39) Acenaphthene-d10	14.32	164	713235	20.00	ng	0.00
64) Phenanthrene-d10	17.08	188	962682	20.00	ng	0.00
76) Chrysene-d12	21.17	240	1372722	20.00	ng	0.00
86) Perylene-d12	23.41	264	1749273	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	1674896	74.85	ng	0.00
7) Phenol-d6	6.86	99	2025391	65.20	ng	0.00
23) Nitrobenzene-d5	8.82	82	1194593	46.31	ng	-0.01
42) 2,4,6-Tribromophenol	15.82	330	767285	86.56	ng	0.00
45) 2-Fluorobiphenyl	12.94	172	2817022	53.96	ng	-0.01
79) Terphenyl-d14	19.66	244	2896166	42.24	ng	0.00
Target Compounds						
10) Phenol	6.89	94	67205	2.148	ng	94
32) Benzoic acid	9.74	122	11628	8.438	ng	84
50) Dimethylphthalate	13.78	163	423312	7.588	ng	# 94
75) Fluoranthene	19.10	202	242624	3.730	ng	78
78) Pyrene	19.45	202	254261	2.650	ng	# 86
88) Benzo(b)fluoranthene	22.73	252	329791m	2.672	ng	

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

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