

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071520\
 Data File : BP002761.D
 Acq On : 15 Jul 2020 19:03
 Operator : CG/JU
 Sample : L3304-03
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 FK-GF-02-058-B

Manual Integrations
 APPROVED

mohammad
 7/16/2020 11:07:23 AM

Quant Time: Jul 16 01:01:34 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 13 10:54:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	174475	20.00	ng	0.00
21) Naphthalene-d8	10.33	136	687664	20.00	ng	0.00
39) Acenaphthene-d10	14.21	164	396490	20.00	ng	0.00
64) Phenanthrene-d10	16.97	188	777355	20.00	ng	0.00
76) Chrysene-d12	21.08	240	619176	20.00	ng	0.00
86) Perylene-d12	23.27	264	565765	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.19	112	770825	74.54	ng	0.00
7) Phenol-d6	6.76	99	1045848	70.86	ng	0.00
23) Nitrobenzene-d5	8.70	82	644068	50.09	ng	0.00
42) 2,4,6-Tribromophenol	15.72	330	263113	63.57	ng	0.00
45) 2-Fluorobiphenyl	12.82	172	1251884	48.84	ng	0.00
79) Terphenyl-d14	19.56	244	1395485	51.67	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.67	163	205930	6.810	ng	99
71) Phenanthrene	17.01	178	514230	12.218	ng	99
72) Anthracene	17.10	178	156049	3.778	ng	98
75) Fluoranthene	19.00	202	785501	16.107	ng	99
78) Pyrene	19.35	202	795058	18.598	ng	98
81) Benzo(a)anthracene	21.06	228	377263	9.413	ng	99
83) Chrysene	21.12	228	349644	9.164	ng	97
87) Indeno(1,2,3-cd)pyrene	25.49	276	189875	4.643	ng	96
88) Benzo(b)fluoranthene	22.62	252	373891m	10.643	ng	
89) Benzo(k)fluoranthene	22.66	252	133644m	3.862	ng	
90) Benzo(a)pyrene	23.17	252	301800	9.066	ng	# 89
92) Benzo(a,h,i)perylene	26.16	276	200150	5.986	ng	96

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

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