

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081720\
 Data File : BP003010.D
 Acq On : 17 Aug 2020 19:27
 Operator : CG/JU
 Sample : L3691-15
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 S708-N-SO-0-0.5-081320

Manual Integrations
 APPROVED

mohammad
 8/19/2020 2:20:53 PM

Quant Time: Aug 18 03:01:12 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	7.69	152	392635	20.00	ng	0.00
21) Naphthalene-d8	10.46	136	1317123	20.00	ng	0.00
39) Acenaphthene-d10	14.32	164	732473	20.00	ng	0.00
64) Phenanthrene-d10	17.07	188	1699599	20.00	ng	0.00
76) Chrysene-d12	21.17	240	1665032	20.00	ng	0.00
86) Perylene-d12	23.42	264	1802846	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	1702275	66.35	ng	0.00
7) Phenol-d6	6.86	99	1994632	56.00	ng	0.00
23) Nitrobenzene-d5	8.83	82	1162876	47.82	ng	0.00
42) 2,4,6-Tribromophenol	15.82	330	680994	74.80	ng	0.00
45) 2-Fluorobiphenyl	12.95	172	2401619	44.80	ng	0.00
79) Terphenyl-d14	19.65	244	2831996	34.05	ng	0.00
Target Compounds						
10) Phenol	6.89	94	81202	2.264	ng	94
50) Dimethylphthalate	13.78	163	540274	9.431	ng	100
71) Phenanthrene	17.12	178	1064320	10.679	ng	99
75) Fluoranthene	19.09	202	1880191	16.374	ng	98
78) Pyrene	19.45	202	1590243	13.662	ng	98
80) Butylbenzylphthalate	20.33	149	5275	3.162	ng	96
81) Benzo(a)anthracene	21.15	228	833933	7.353	ng	99
83) Chrysene	21.20	228	867049	8.104	ng	97
87) Indeno(1,2,3-cd)pyrene	25.70	276	619859	4.293	ng	95
88) Benzo(b)fluoranthene	22.74	252	1353886m	10.642	ng	
89) Benzo(k)fluoranthene	22.78	252	442019m	3.676	ng	
90) Benzo(a)pyrene	23.32	252	844698	7.348	ng	99
92) Benzo(a,h,i)perylene	26.39	276	624575	5.256	ng	100

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

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