

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081720\
 Data File : BP003008.D
 Acq On : 17 Aug 2020 18:19
 Operator : CG/JU
 Sample : L3691-14
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 S707-N-SO-0-0.5-081320

Manual Integrations
 APPROVED

mohammad
 8/20/2020 12:20:29 PM

Quant Time: Aug 18 05:44:57 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	361360	20.00	ng	0.00
21) Naphthalene-d8	10.46	136	1271409	20.00	ng	0.00
39) Acenaphthene-d10	14.32	164	738245	20.00	ng	0.00
64) Phenanthrene-d10	17.07	188	1545996	20.00	ng	0.00
76) Chrysene-d12	21.17	240	1685589	20.00	ng	0.00
86) Perylene-d12	23.42	264	1889860	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	1843104	78.06	ng	0.00
7) Phenol-d6	6.86	99	2189257	66.79	ng	0.00
23) Nitrobenzene-d5	8.83	82	1115223	47.51	ng	0.00
42) 2,4,6-Tribromophenol	15.82	330	776845	84.66	ng	0.00
45) 2-Fluorobiphenyl	12.95	172	2796193	51.75	ng	0.00
79) Terphenyl-d14	19.65	244	3856458	45.80	ng	0.00
Target Compounds						
10) Phenol	6.89	94	89462	2.710	ng	93
32) Benzoic acid	9.75	122	5227m	8.058	ng	
50) Dimethylphthalate	13.78	163	591591	10.246	ng	99
71) Phenanthrene	17.12	178	251779	2.777	ng	98
75) Fluoranthene	19.09	202	701351	6.715	ng	100
78) Pyrene	19.44	202	614056	5.211	ng	97
81) Benzo(a)anthracene	21.15	228	285404	2.486	ng	99
83) Chrysene	21.20	228	309064	2.854	ng	98
88) Benzo(b)fluoranthene	22.74	252	466556	3.498	ng	98
90) Benzo(a)pyrene	23.32	252	310191	2.574	ng	98

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

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