

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP072420\
 Data File : BP002865.D
 Acq On : 24 Jul 2020 21:39
 Operator : CG/JU
 Sample : L3382-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CTR-007

Manual Integrations
 APPROVED

mohammad
 7/27/2020 12:22:15 PM

Quant Time: Jul 25 00:27:12 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 24 15:50:46 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.53	152	38333	20.00	ng	0.00
21) Naphthalene-d8	10.31	136	159373	20.00	ng	0.00
39) Acenaphthene-d10	14.19	164	105770	20.00	ng	0.00
64) Phenanthrene-d10	16.95	188	263556	20.00	ng	0.00
76) Chrysene-d12	21.05	240	365739	20.00	ng	0.00
86) Perylene-d12	23.23	264	439240	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.16	112	213328	93.89	ng	0.00
7) Phenol-d6	6.74	99	295943	91.27	ng	0.00
23) Nitrobenzene-d5	8.69	82	169156	56.77	ng	0.00
42) 2,4,6-Tribromophenol	15.70	330	105797	95.82	ng	0.00
45) 2-Fluorobiphenyl	12.80	172	399777	58.47	ng	0.00
79) Terphenyl-d14	19.53	244	789602	49.50	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	13.66	163	64577	8.006	ng	99
71) Phenanthrene	16.99	178	40514	2.839	ng	97
75) Fluoranthene	18.98	202	66980	4.051	ng	98
78) Pyrene	19.33	202	89451	3.542	ng	98
81) Benzo(a)anthracene	21.04	228	66378	2.804	ng	97
83) Chrysene	21.09	228	99684	4.423	ng	96
87) Indeno(1,2,3-cd)pyrene	25.43	276	86521	2.725	ng	99
88) Benzo(b)fluoranthene	22.59	252	195095	7.153	ng	98
89) Benzo(k)fluoranthene	22.62	252	57231m	2.130	ng	
90) Benzo(a)pyrene	23.14	252	92713	3.587	ng	98
92) Benzo(g,h,i)perylene	26.10	276	84271	3.246	ng	97

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

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