

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP010620\
 Data File : BP001531.D
 Acq On : 06 Jan 2020 15:58
 Operator : JU
 Sample : L1028-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 NORTH-BERGEN

Manual Integrations
 APPROVED

mohammad
 1/7/2020 12:31:21 PM

Quant Time: Jan 06 16:49:33 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP010320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 06 16:38:29 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	47616	20.00	ng	0.00
21) Naphthalene-d8	10.46	136	177459	20.00	ng	0.00
39) Acenaphthene-d10	14.32	164	122836	20.00	ng	0.00
64) Phenanthrene-d10	17.08	188	262314	20.00	ng	0.00
76) Chrysene-d12	21.18	240	236382	20.00	ng	0.00
87) Perylene-d12	23.43	264	241371	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	233283	82.31	ng	0.00
7) Phenol-d6	6.89	99	339573	88.95	ng	0.00
23) Nitrobenzene-d5	8.83	82	237305	66.05	ng	0.00
42) 2,4,6-Tribromophenol	15.82	330	151190	108.75	ng	0.00
45) 2-Fluorobiphenyl	12.94	172	449326	53.79	ng	0.00
79) Terphenyl-d14	19.66	244	647038	54.55	ng	0.00
Target Compounds						
10) Phenol	6.91	94	9454	2.354	ng	85
50) Dimethylphthalate	13.79	163	28057	2.922	ng	# 99
71) Phenanthrene	17.12	178	151988	10.683	ng	98
72) Anthracene	17.21	178	41272	2.927	ng	99
75) Fluoranthene	19.10	202	326269	18.944	ng	99
78) Pyrene	19.45	202	294804	17.187	ng	98
81) Benzo(a)anthracene	21.16	228	172564	10.404	ng	98
83) Chrysene	21.22	228	155880	9.802	ng	97
86) Indeno(1,2,3-cd)pyrene	25.70	276	82108	4.242	ng	98
88) Benzo(b)fluoranthene	22.76	252	183956	12.309	ng	# 94
89) Benzo(k)fluoranthene	22.79	252	52872m	3.693	ng	
90) Benzo(a)pyrene	23.33	252	124762	8.918	ng	97
92) Benzo(a,h,i)perylene	26.39	276	78616	5.669	ng	98

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

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