

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP010620\
 Data File : BP001538.D
 Acq On : 06 Jan 2020 19:54
 Operator : JU
 Sample : K6465-03
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GP-10-(1-3)

Manual Integrations
 APPROVED

mohammad
 1/7/2020 12:31:34 PM

Quant Time: Jan 07 03:20:37 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	48097	20.00	ng	0.00
21) Naphthalene-d8	10.46	136	179332	20.00	ng	0.00
39) Acenaphthene-d10	14.32	164	120864	20.00	ng	0.00
64) Phenanthrene-d10	17.08	188	270688	20.00	ng	0.00
76) Chrysene-d12	21.18	240	216002	20.00	ng	0.00
87) Perylene-d12	23.43	264	229078	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	249764	87.24	ng	0.00
7) Phenol-d6	6.89	99	366371	95.01	ng	0.00
23) Nitrobenzene-d5	8.83	82	252938	69.66	ng	0.00
42) 2,4,6-Tribromophenol	15.82	330	157432	115.09	ng	0.00
45) 2-Fluorobiphenyl	12.94	172	542992	66.06	ng	0.00
79) Terphenyl-d14	19.66	244	768799	70.93	ng	0.00
Target Compounds						
10) Phenol	6.91	94	13719	3.382	ng	95
50) Dimethylphthalate	13.79	163	48506	5.134	ng	99
71) Phenanthrene	17.12	178	298071	20.302	ng	99
72) Anthracene	17.21	178	74779	5.139	ng	98
75) Fluoranthene	19.10	202	569846	32.063	ng	99
78) Pyrene	19.45	202	536990	34.259	ng	99
81) Benzo(a)anthracene	21.16	228	262204	17.300	ng	98
83) Chrysene	21.22	228	230481	15.860	ng	98
86) Indeno(1,2,3-cd)pyrene	25.70	276	128424	7.261	ng	98
88) Benzo(b)fluoranthene	22.76	252	274859	19.378	ng	# 94
89) Benzo(k)fluoranthene	22.80	252	78856m	5.804	ng	
90) Benzo(a)pyrene	23.33	252	201955	15.211	ng	97
91) Dibenzo(a,h)anthracene	25.71	278	34699	2.630	ng	# 89
92) Benzo(a,h,i)perylene	26.39	276	124009	9.422	ng	97

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

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