

Data Path : Z:\svoasrv\HPCHEM1\BNA P\Data\BP100419\
 Data File : BP000512.D
 Acq On : 04 Oct 2019 17:12
 Operator : JU
 Sample : K5160-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CSA-3-2

Manual Integrations
 APPROVED

mohammad
 10/7/2019 4:38:16 PM

Quant Time: Oct 23 05:17:08 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP100119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 01 18:03:42 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.76	152	251018	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	918450	20.00	ng	-0.01
39) Acenaphthene-d10	9.81	164	517266	20.00	ng	0.00
64) Phenanthrene-d10	11.31	188	972011	20.00	ng	0.00
76) Chrysene-d12	13.98	240	848299	20.00	ng	0.00
87) Perylene-d12	15.46	264	916726	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	1413721	90.53	ng	0.00
7) Phenol-d6	6.40	99	1796783	95.48	ng	0.00
23) Nitrobenzene-d5	7.32	82	1006004	66.90	ng	-0.01
42) 2,4,6-Tribromophenol	10.60	330	568630	103.16	ng	0.00
45) 2-Fluorobiphenyl	9.13	172	2113802	66.12	ng	0.00
79) Terphenyl-d14	12.92	244	2641839	63.06	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.52	163	516746	14.377	ng	99
75) Fluoranthene	12.53	202	242486	4.423	ng	98
78) Pyrene	12.76	202	241047	4.120	ng	100
81) Benzo(a)anthracene	13.97	228	167604m	3.238	ng	
83) Chrysene	14.00	228	161941	3.188	ng	98
88) Benzo(b)fluoranthene	15.03	252	258037m	4.961	ng	
90) Benzo(a)pyrene	15.39	252	179289	3.698	ng	96
92) Benzo(g,h,i)perylene	17.36	276	130698	2.735	ng	100

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

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