

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120219\
 Data File : BP001130.D
 Acq On : 02 Dec 2019 17:14
 Operator : JU
 Sample : K6022-07
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-16-S

Manual Integrations
 APPROVED

mohammad
 12/4/2019 4:59:15 PM

Quant Time: Dec 02 17:54:41 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP112719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 27 17:38:46 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.33	152	166731	20.00	ng	0.00
21) Naphthalene-d8	10.36	136	638179	20.00	ng	0.00
39) Acenaphthene-d10	13.23	164	342314	20.00	ng	0.00
64) Phenanthrene-d10	15.62	188	590938	20.00	ng	0.00
76) Chrysene-d12	19.27	240	423756	20.00	ng	0.00
87) Perylene-d12	21.05	264	463403	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	6.23	112	1035209	103.01	ng	0.00
7) Phenol-d6	7.77	99	1452177	108.46	ng	0.00
23) Nitrobenzene-d5	9.21	82	951897	64.71	ng	0.00
42) 2,4,6-Tribromophenol	14.52	330	286890	87.44	ng	0.00
45) 2-Fluorobiphenyl	12.15	172	1473724	63.01	ng	0.00
79) Terphenyl-d14	17.91	244	1388271	60.61	ng	0.00
Target Compounds						
50) Dimethylphthalate	12.81	163	253709	9.910	ng	100
71) Phenanthrene	15.66	178	150309	4.838	ng	99
75) Fluoranthene	17.37	202	221660	6.739	ng	99
78) Pyrene	17.67	202	197053	5.904	ng	97
81) Benzo(a)anthracene	19.26	228	96328	3.279	ng	93
83) Chrysene	19.30	228	91489	3.477	ng	97
88) Benzo(b)fluoranthene	20.56	252	102538m	3.623	ng	
90) Benzo(a)pyrene	20.97	252	79888	3.064	ng	99

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

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