

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071620\
 Data File : BP002782.D
 Acq On : 16 Jul 2020 20:45
 Operator : CG/JU
 Sample : L3184-07 5X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EO-01-071520

Manual Integrations
 APPROVED

mohammad
 7/17/2020 12:30:43 PM

Quant Time: Jul 17 01:19:42 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 13 10:54:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	249715	20.00	ng	0.01
21) Naphthalene-d8	10.33	136	982842	20.00	ng	0.00
39) Acenaphthene-d10	14.21	164	535845	20.00	ng	0.00
64) Phenanthrene-d10	16.97	188	876908	20.00	ng	0.00
76) Chrysene-d12	21.08	240	611248	20.00	ng	0.00
86) Perylene-d12	23.27	264	629162	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.19	112	225441	15.23	ng	0.00
7) Phenol-d6	6.76	99	309501	14.65	ng	0.00
23) Nitrobenzene-d5	8.70	82	199028	10.83	ng	0.00
42) 2,4,6-Tribromophenol	15.72	330	72878	13.03	ng	0.00
45) 2-Fluorobiphenyl	12.82	172	422328	12.19	ng	0.00
79) Terphenyl-d14	19.56	244	306691	11.50	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.67	163	82982	2.031	ng	99
71) Phenanthrene	17.01	178	250196	5.270	ng	100
75) Fluoranthene	19.00	202	332491	6.044	ng	97
78) Pyrene	19.35	202	311548	7.382	ng	99
81) Benzo(a)anthracene	21.06	228	131074	3.313	ng	97
83) Chrysene	21.12	228	119027m	3.160	ng	
88) Benzo(b)fluoranthene	22.62	252	130063	3.329	ng	# 89
90) Benzo(a)pyrene	23.17	252	97882	2.644	ng	# 90

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

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