

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072219\
 Data File : BF115713.D
 Acq On : 22 Jul 2019 17:00
 Operator : HP/JU
 Sample : K3928-01 2X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ROLL-OFF-COMP

Manual Integrations
 APPROVED

mohammad
 7/23/2019 4:24:33 PM

Quant Time: Jul 23 03:52:26 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 12 14:03:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.88	152	169304	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	652989	20.00	ng	0.00
39) Acenaphthene-d10	9.92	164	312937	20.00	ng	-0.01
64) Phenanthrene-d10	11.40	188	474843	20.00	ng	0.00
76) Chrysene-d12	14.04	240	404860	20.00	ng	0.00
87) Perylene-d12	15.53	264	366714	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	357729	34.56	ng	-0.01
7) Phenol-d6	6.50	99	468230	35.65	ng	-0.01
23) Nitrobenzene-d5	7.43	82	279783	24.91	ng	-0.02
42) 2,4,6-Tribromophenol	10.70	330	102220	27.98	ng	-0.01
45) 2-Fluorobiphenyl	9.23	172	537287	30.05	ng	-0.02
79) Terphenyl-d14	12.99	244	474734	21.64	ng	-0.01
Target Compounds						
50) Dimethylphthalate	9.63	163	48059	2.041	ng	Qvalue 98
71) Phenanthrene	11.43	178	52928	2.175	ng	96
75) Fluoranthene	12.62	202	147882	5.749	ng	98
78) Pyrene	12.84	202	143680	4.189	ng	97
81) Benzo(a)anthracene	14.03	228	84632	3.173	ng	97
83) Chrysene	14.07	228	76706	2.865	ng	96
88) Benzo(b)fluoranthene	15.09	252	97487m	4.128	ng	
90) Benzo(a)pyrene	15.46	252	58466	2.881	ng	# 85

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

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