

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052419\
 Data File : BM020566.D
 Acq On : 25 May 2019 06:20
 Operator : JU/SJ
 Sample : K3026-03 5X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TP-5-S

Manual Integrations
 APPROVED

mohammad
 5/28/2019 8:41:34 AM

Quant Time: May 27 03:53:47 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM052419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 24 16:07:05 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	94454	20.00	ng	0.00
21) Naphthalene-d8	10.45	136	366625	20.00	ng	0.00
39) Acenaphthene-d10	14.32	164	259152	20.00	ng	0.00
64) Phenanthrene-d10	17.07	188	605329	20.00	ng	0.00
76) Chrysene-d12	21.27	240	458186	20.00	ng	0.00
87) Perylene-d12	23.51	264	517523	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.25	112	104376	21.68	ng	0.00
7) Phenol-d6	6.84	99	147410	20.29	ng	0.00
23) Nitrobenzene-d5	8.83	82	96554	11.05	ng	0.00
42) 2,4,6-Tribromophenol	15.82	330	92801	17.87	ng	0.00
45) 2-Fluorobiphenyl	12.93	172	178117	8.62	ng	0.00
79) Terphenyl-d14	19.70	244	261621	10.32	ng	0.00
Target Compounds						
71) Phenanthrene	17.12	178	71477	2.174	ng	99
75) Fluoranthene	19.14	202	133045	2.909	ng	98
78) Pyrene	19.50	202	117515	4.085	ng	97
81) Benzo(a)anthracene	21.25	228	64244	2.026	ng	# 93
83) Chrysene	21.30	228	63920	2.086	ng	# 93
88) Benzo(b)fluoranthene	22.83	252	80651m	2.381	ng	

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

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